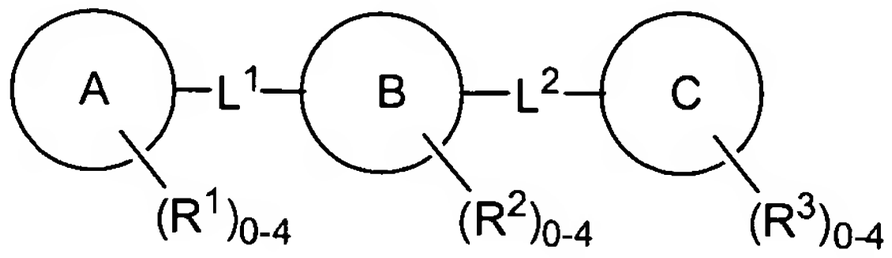
Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. (original) A compound for modulating c-Kit activity according to Formula **I**,



I

or a pharmaceutically acceptable salt, hydrate, or prodrug thereof, wherein,

ring A is a five- to fourteen-membered heteroaryl;

each R^1 is independently selected from -H, halogen, trihalomethyl, -CN, $-NO_2$, $-OR^4$, $-N(R^4)R^4$, $-S(O)_{0-2}R^4$, $-SO_2N(R^4)R^4$, $-CO_2R^4$, $-C(=O)N(R^4)R^4$, $-C(=NR^5)N(R^4)R^4$, $-C(=NR^5)R^4$, $-N(R^4)SO_2R^4$, $-N(R^4)C(O)R^4$, $-NCO_2R^4$, $-C(=O)R^4$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

two adjacent of R^1, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R^{10};

L^1 is selected from a single bond, an optionally substituted C_{1-2}alkylene, -O-, $-CH_2O-$, $-N(R^7)-$, $-C(=O)N(R^7)-$, $-SO_2N(R^7)-$, $-CH_2N(R^7)-$, and $-S(O)_{0-2}-$;

ring B is a five- to ten-membered aryl or a five- to ten-membered heterocyclyl;

each R^2 is independently selected from -H, halogen, trihalomethyl, -CN, $-NO_2$, $-OR^4$, $-N(R^4)R^4$, $-S(O)_{0-2}R^4$, $-SO_2N(R^4)R^4$, $-CO_2R^4$, $-C(=O)N(R^4)R^4$, $-C(=NR^5)N(R^4)R^4$, $-C(=NR^5)R^4$, $-N(R^4)SO_2R^4$, $-N(R^4)C(O)R^4$, $-NCO_2R^4$, $-C(=O)R^4$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-}

$_6$alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_1-$_6$alkyl;

two adjacent of R^2, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R^{15};

L^2 is a selected from C_4alkylene, C_4alkylidene, C_4alkylidyne, $-X(CH_2)_2O-$, $-X(CH_2)_2N(R^7)-$, $-XCH_2SO_2N(R^7)-$, $-XN(R^7)C(=O)N(R^7)-$, $-XCH_2C(=O)N(R^7)-$, $-(CH_2)_3X-$, $-XN(R^7)SO_2N(R^7)-$, $-XCH_2N(R^7)SO_2-$, $-CH_2X(CH_2)_2-$, $-CH=CHC(=O)N(R^7)-$, $-CH=CHSO_2N(R^7)-$, $-XCH_2N(R^7)C(=O)-$, -M-M-, $-CH_2N(R^7)C(=O)O-$, and $-CH_2OC(=O)N(R^7)-$; wherein X is selected from $-CH_2-$, -O-, $-N(R^7)-$, -C(=O)-, and $-S(O)_{0-2}-$; M is selected from $-C(=O)N(R^7)-$ and $-SO_2N(R^7)-$; and any C-H of L^2 is optionally C-R^{20};

ring C is either a five- to ten-membered aryl or a five- to ten-membered heteroaryl;

each R^3 is independently selected from -H, halogen, trihalomethyl, -CN, $-NO_2$, $-OR^4$, $-N(R^4)R^4$, $-S(O)_{0-2}R^4$, $-SO_2N(R^4)R^4$, $-CO_2R^4$, $-C(=O)N(R^4)R^4$, $-C(=NR^5)N(R^4)R^4$, $-C(=NR^5)R^4$, $-N(R^4)SO_2R^4$, $-N(R^4)C(O)R^4$, $-NCO_2R^4$, $-C(=O)R^4$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_1-$_6$alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_1-$_6$alkyl; provided R^3 is not a cyclic sulfonamide attached to ring C via the nitrogen of said cyclic sulfonamide;

two adjacent of R^3, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R^{25};

R^4 is selected from -H, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

two of R^4, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted

five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R^5 is selected from -H, -CN, $-NO_2$, $-OR^4$, $-S(O)_{0-2}R^4$, $-CO_2R^4$, optionally substituted C_{1-6}alkyl, optionally substituted C_{1-6}alkenyl, and optionally substituted C_{1-6}alkynyl;

R^7 is selected from -H, optionally substituted C_{1-6}alkyl, $-SO_2N(R^4)R^4$, $-CO_2R^4$, $-C(=O)N(R^4)R^4$, $-C(=NR^5)N(R^4)R^4$, $-C(=NR^5)R^4$, $-C(=O)R^4$, optionally substituted alkoxy, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl; and

each of R^{10}, each of R^{15}, each of R^{20}, and each of R^{25} is independently selected from -H, halogen, trihalomethyl, -CN, $-NO_2$, $-OR^4$, $-N(R^4)R^4$, $-S(O)_{0-2}R^4$, $-SO_2N(R^4)R^4$, $-CO_2R^4$, $-C(=O)N(R^4)R^4$, $-C(=NR^5)N(R^4)R^4$, $-C(=NR^5)R^4$, $-N(R^4)SO_2R^4$, $-N(R^4)C(O)R^4$, $-NCO_2R^4$, $-C(=O)R^4$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

provided:

1) when both ring B and ring C are phenyl:

a) and the compound comprises ring B-$CH_2N(H)C(=O)N(H)$-ring C, then L^1 must be a single bond; R^3 can not comprise a group of the formula $-O(CH_2)_{2-4}$-*N*-piperazine that is *ortho*- to L^2; and ring A cannot be a 5-methyl-[1,2,4]-oxadiazol-3-yl radical, a 4H-[1,2,4]-oxadiazol-5-one-3-yl radical, nor a 4'-[2,2';6',2"]terpyridinyl radical;

b) and L^1 is single bond, then L^2 cannot comprise -N(H)C(=O)C(=O)N(H)- nor -N(H)C(=Q)C(H)CNC(=O)- (where Q is S or O);

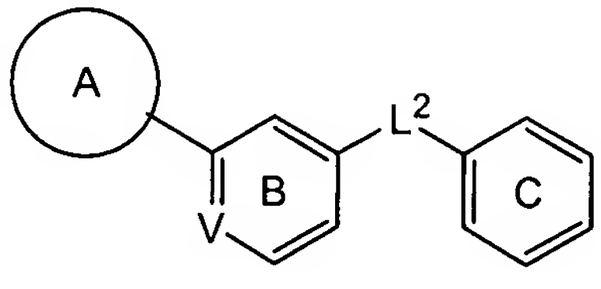
c) and L^1 is other than single bond, then A cannot be quinolin-2-yl-L^1, quinolin-3-yl-L^1, or quinolin-4-yl-L^1;

2) when ring A is a fused aryl system, then L^1 must be a single bond;

3) when ring B is phenyl, ring C is a C_{6-16}carbocyclic, L^1 is a single bond, and the compound comprises –ring B-$OCH_2C(=O)N(H)$- then ring A cannot be a 2,5-dimethyl-1H-pyrrole-1-yl radical;

4) ring A cannot be a pyrimidin-2-yl radical when L^1 is -N(H)- and ring B is phenyl;

5) when the compound comprises the formula,



where V is =C(H)- or =N-, and there is a nitrogen of L^2 bound directly to ring B, then A can not comprise a [1,2,4]-oxadiazol-3-yl radical; and

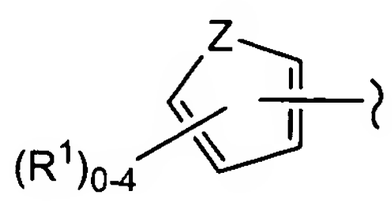
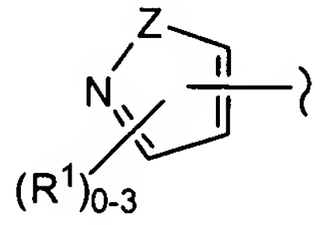
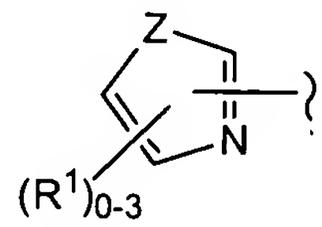
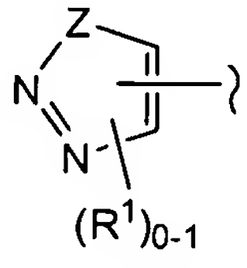
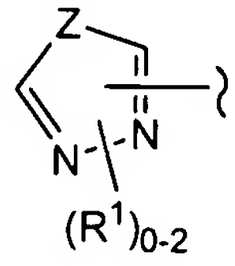
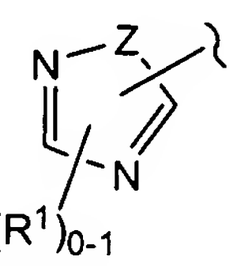
6) the compound is not one of: N-naphthalen-1-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(phenyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,6-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,6-diethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(ethyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(dimethylamino)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-chloro-3-methylphenyl)-2-{[3-(1H-tetrazol-1-

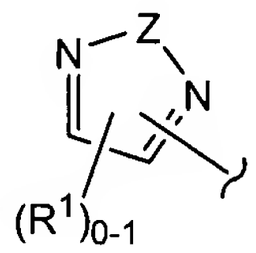
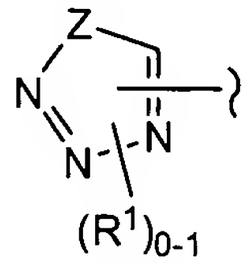
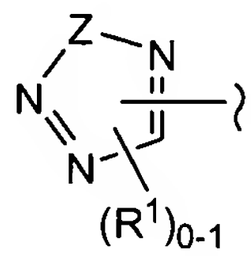
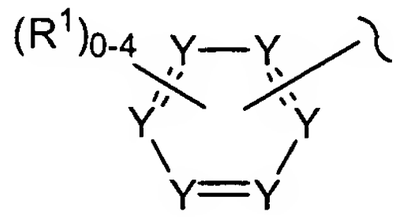
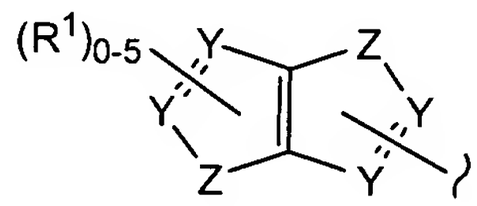
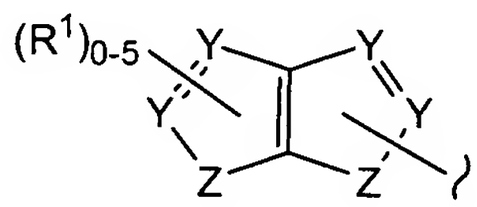
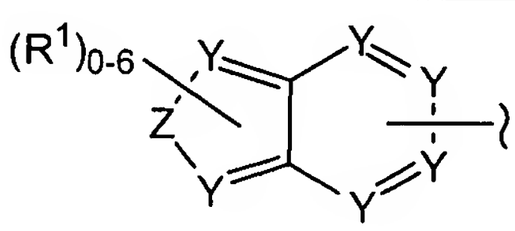
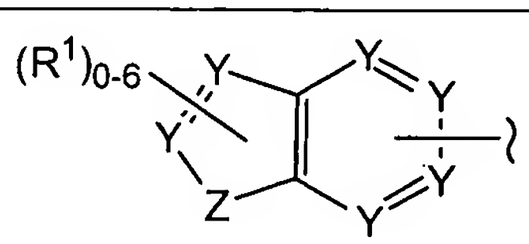
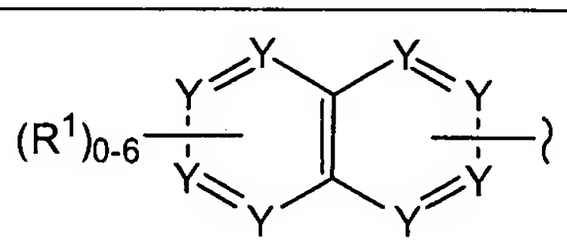
yl)phenyl]oxy}acetamide, N-(4-bromophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoromethyl)phenyl]acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl] oxy}-N-[3-(trifluoromethyl)phenyl]acetamide, methyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoate, ethyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoate, 3-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoic acid, N-[3-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-chloro-5-(trifluoromethyl) phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acet-amide, N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-aminophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, and N-(4-acetylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide.

2. (original) The compound according to claim 1, wherein L^1 is a single bond.

3. (original) The compound according to claim 2, wherein ring A contains between one and four annular nitrogens.

4. (original) The compound according to claim 3, wherein ring A is selected from the following:

wherein each Y is independently either =C(H)- or =N-; and Z is selected from -O-, -$S(O)_{0-2}$-, and -$N(R^7)$-.

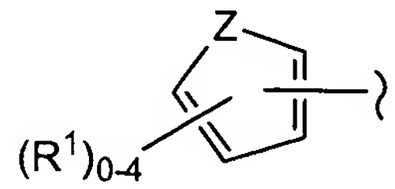
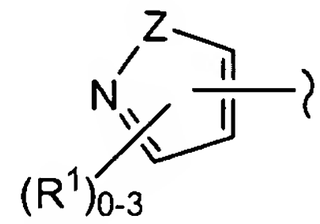
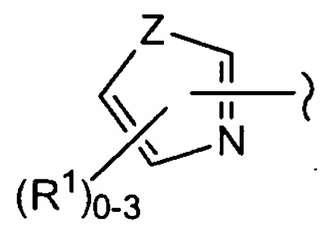
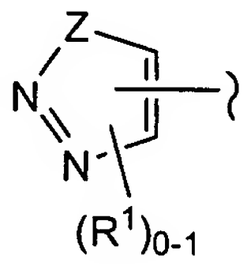
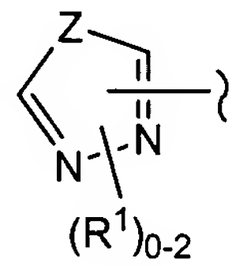
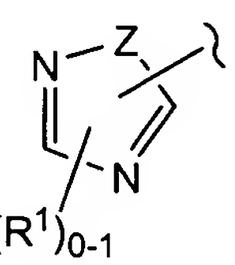
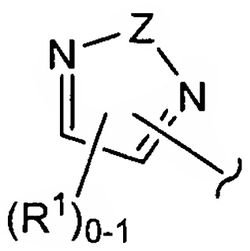
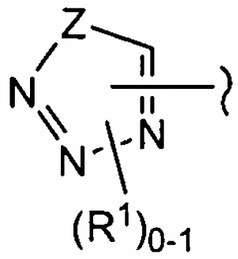
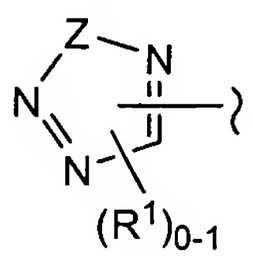
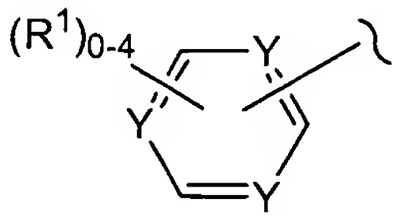
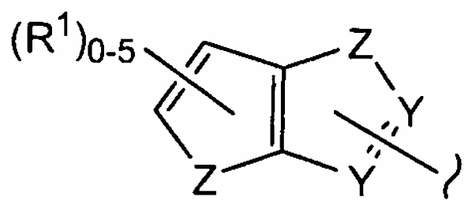
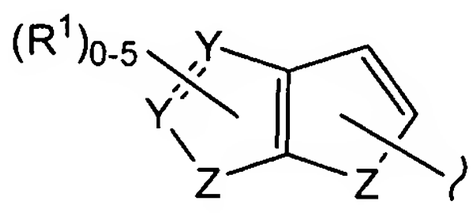
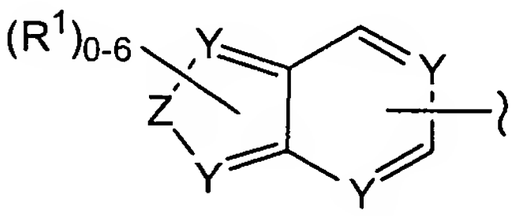
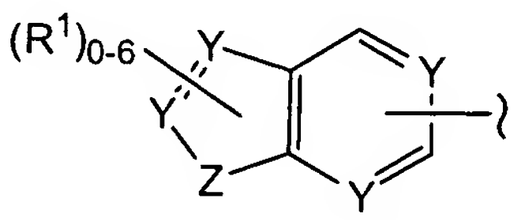
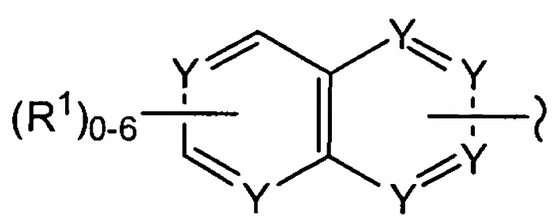
5. (original) The compound according to claim 4, wherein ring B is phenylene or pyridylene.

6. (original) The compound according to claim 5, wherein the annular atoms of ring B to which L^1 and L^2 are attached are not contiguous.

7. (original) The compound according to claim 6, wherein L^2 is selected from -$X(CH_2)_2O$-, -$X(CH_2)_2N(R^7)$-, -$CH_2XC(=O)N(R^7)$-, -$XCH_2SO_2N(R^7)$-, -$XN(R^7)C(=O)N(R^7)$- and -$XCH_2C(=O)N(R^7)$-; wherein X is selected from -CH_2-, -O-, -$S(O)_{0-2}$- and -$N(R^7)$-; and any C-H of L^2 is optionally C-R^{20}.

8. (original) The compound according to claim 7, wherein L^2 is selected from -$N(H)N(H)C(=O)N(H)$-, -$CH_2N(H)C(=O)N(H)$-, -$CH_2OC(=O)N(H)$-, and -$XCH_2C(=O)N(H)$-; wherein X is selected from -O-, -$S(O)_{0-2}$-, and -$N(R^7)$-; and any C-H of L^2 is optionally C-R^{20}.

9. (original) The compound according to claim 8, wherein ring A is selected from the following:

wherein each Y is independently either =C(H)- or =N-; and Z is selected from -O-, -S-, and -N(R^7)-.

10. (original) The compound according to claim 9, wherein ring C is phenyl or pyridyl.

11. (original) The compound according to claim 10, wherein there exists at least one of R^3 that is halogen.

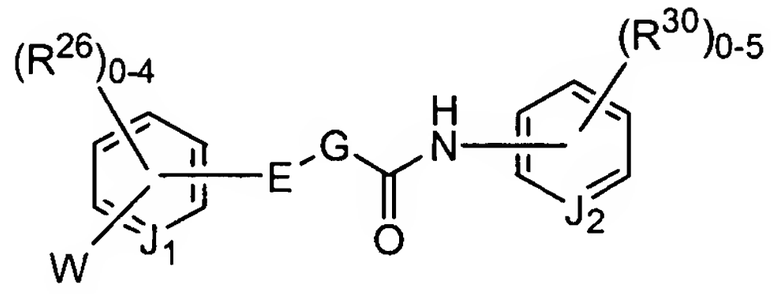
12. (original) The compound according to claim 10, wherein there exists at least one of R^3 that is trihalomethyl.

13. (original) The compound according to claim 10, wherein there exists at least one of R^3 that is trifluoromethyl.

14. (original) The compound according to claim 13, wherein ring C is a phenyl comprising a trifluoromethyl radical *meta*- to L^2.

15. (original) The compound according to claim 10, wherein each of R^3 is independently selected from -H, halogen, trihalomethyl, $-OR^4$, $-CO_2R^4$, $-C(=O)R^4$, and optionally substituted C_{1-6}alkyl.

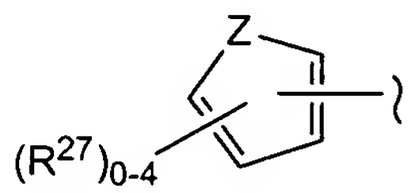
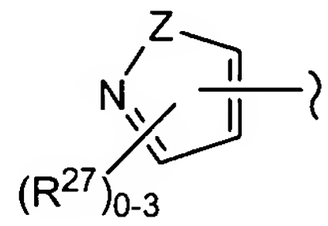
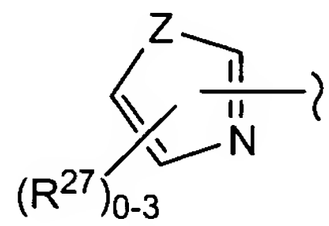
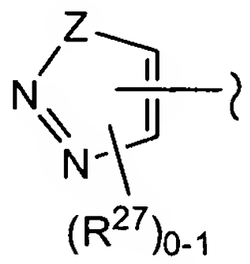
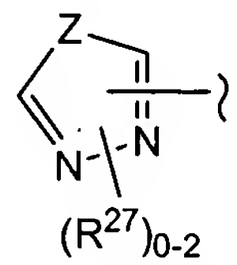
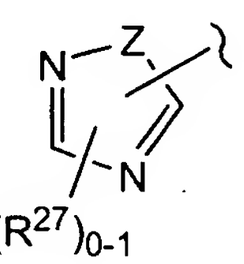
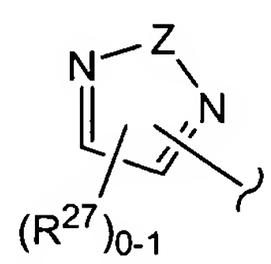
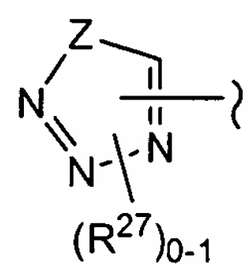
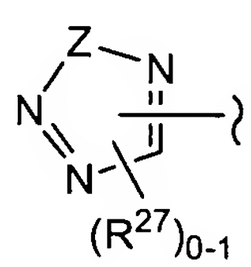
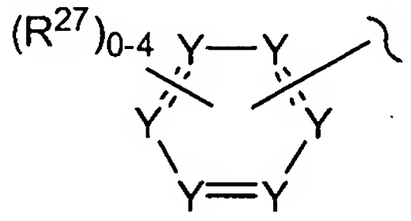
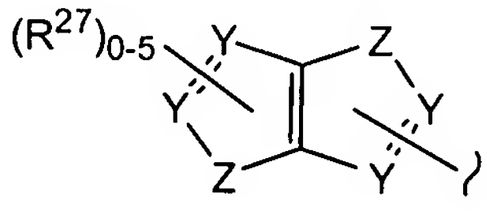
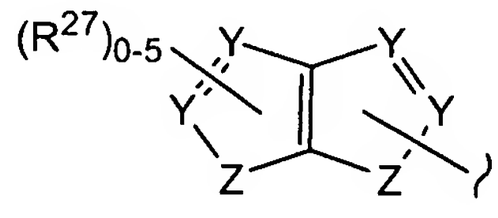
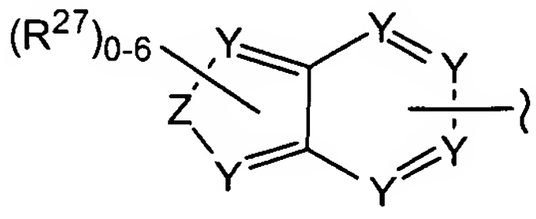
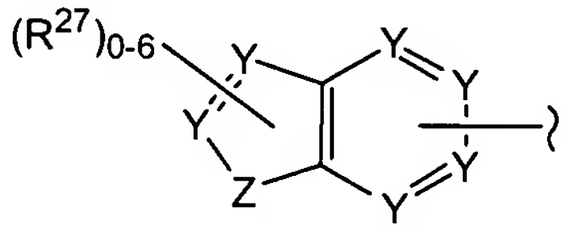
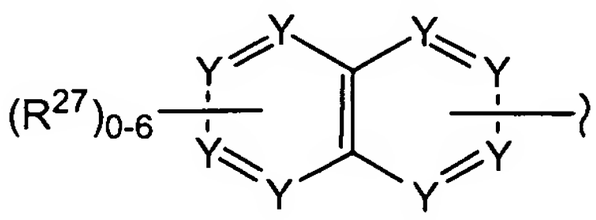
16. (original) A compound for modulating c-Kit activity according to Formula **II**,



II

or a pharmaceutically acceptable salt, hydrate, or prodrug thereof, wherein,

W is selected from the following:

each of R^{27} independently selected from halogen, trihalomethyl, -CN, -NO_2, -OR^{55}, -$N(R^{55})R^{55}$, -$S(O)_{0-2}R^{55}$, -$SO_2N(R^{55})R^{55}$, -CO_2R^{55}, -$C(=O)N(R^{55})R^{55}$, -$C(=NR^{50})N(R^{55})R^{55}$, -$C(=NR^{50})R^{55}$, -$N(R^{55})SO_2R^{55}$, -$N(R^{55})C(O)R^{55}$, -$NCO_2R^{55}$, -$C(=O)R^{55}$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

each Y is independently either =C(H)- or =N-;

Z is selected from -O-, -$S(O)_{0-2}$-, and -$N(R^7)$-

E and G are each independently selected from -O-, -$S(O)_{0-2}$-, -$C(R^{31})R^{32}$-, and -$N(R^{33})$-;

J_1 and J_2 are each independently =C(H)- or =N-;

each of R^{26} and R^{30} is independently selected from -H, halogen, trihalomethyl, -CN, -NO_2, -OR^{40}, -$N(R^{40})R^{40}$, -$S(O)_{0-2}R^{40}$, -$SO_2N(R^{40})R^{40}$, -CO_2R^{40}, -$C(=O)N(R^{40})R^{40}$, -$C(=NR^{50})N(R^{40})R^{40}$, -$C(=NR^{50})R^{40}$, -$N(R^{40})SO_2R^{40}$, -$N(R^{40})C(O)R^{40}$, -$NCO_2R^{40}$, -$C(=O)R^{40}$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

two adjacent of R^{26} or two adjacent of R^{30}, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R^{35};

$R^{31}$and R^{32} are each independently selected from -H, halogen, trihalomethyl, -CN, -NO_2, -OR^{40}, -$N(R^{40})R^{40}$, -$S(O)_{0-2}R^{40}$, -$SO_2N(R^{40})R^{40}$, -CO_2R^{40}, -$C(=O)N(R^{40})R^{40}$, -$C(=NR^{50})N(R^{40})R^{40}$, -$C(=NR^{50})R^{40}$, -$N(R^{40})SO_2R^{40}$, -$N(R^{40})C(O)R^{40}$, -$NCO_2R^{40}$, -$C(=O)R^{40}$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

R^{33} is selected from -H, optionally substituted lower alkyl, -$SO_2N(R^{40})R^{40}$, -CO_2R^{40}, -$C(=O)N(R^{40})R^{40}$, -$C(=NR^{50})N(R^{40})R^{40}$, -$C(=NR^{50})R^{40}$, -$C(=O)R^{40}$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted

aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

R^{40} is selected from -H, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl;

two of R^{40}, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R^{50} is selected from -H, -CN, $-NO_2$, $-OR^{40}$, $-S(O)_{0-2}R^{40}$, $-CO_2R^{40}$, optionally substituted C_{1-6}alkyl, optionally substituted C_{1-6}alkenyl, and optionally substituted C_{1-6}alkynyl;

R^{55} is selected from -H, optionally substituted C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl; and

two of R^{55}, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P.

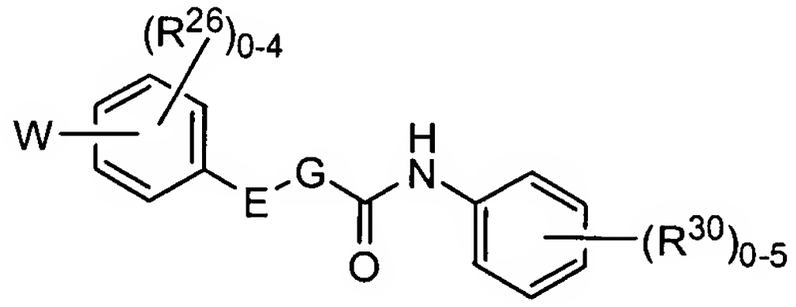
17. (original) The compound according to claim 16, wherein the annular carbons of ring B to which W and E are attached are not contiguous.

18. (original) The compound according to claim 17, wherein R^{30} is selected from -H, halogen, trihalomethyl, $-OR^{40}$, $-N(R^{40})R^{40}$, $-CO_2R^{40}$, $-C(=O)R^{40}$, optionally substituted alkoxy, optionally substituted C_{1-6}alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6}alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6}alkyl.

19. (original) The compound according to claim 18, wherein there exists at least one of R^{30} that is trihalomethyl.

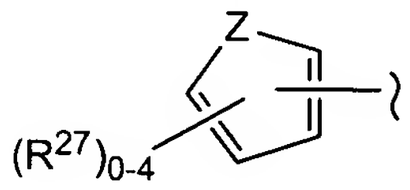
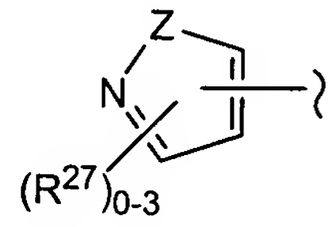
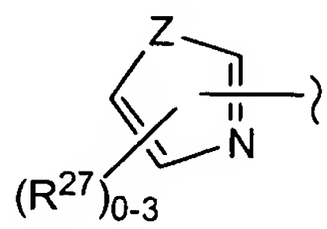
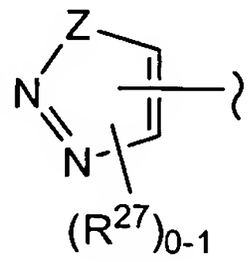
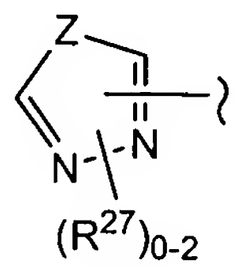
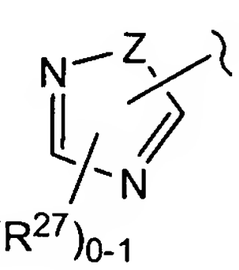
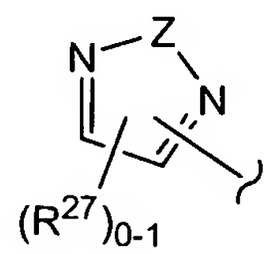
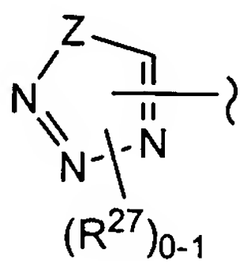
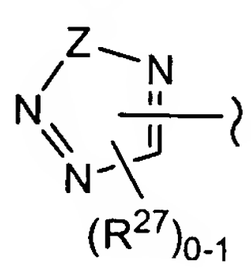
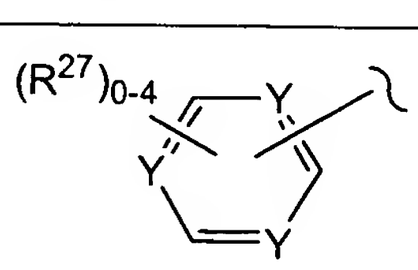
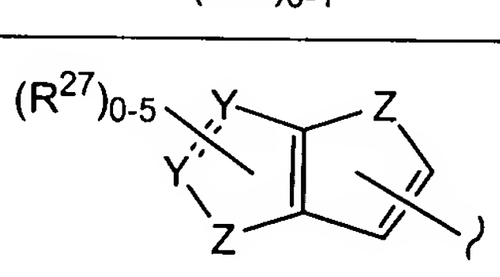
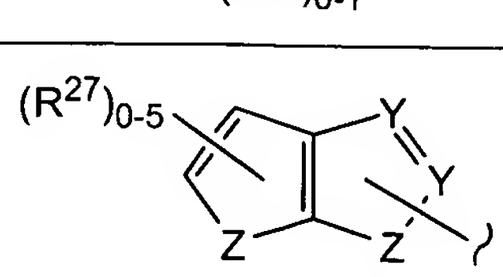
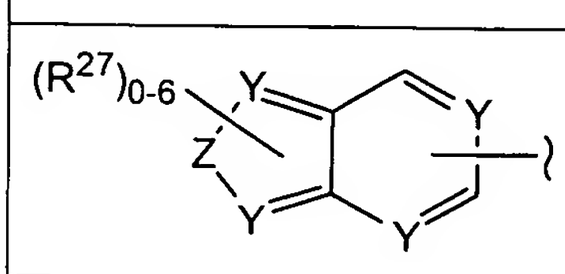
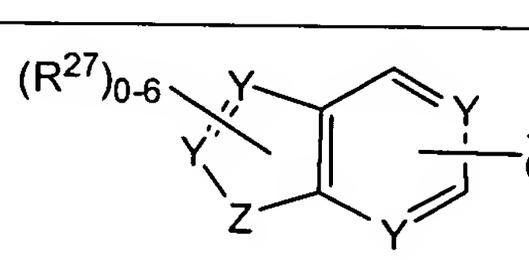
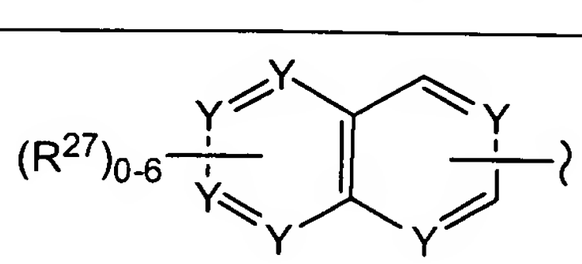
20. (original) The compound according to claim 18, wherein there exists at least one of R^{30} that is trifluoromethyl.

21. (original) The compound according to claim 18, according to formula **III**.



III

22. (original) The compound according to claim 21, wherein W is selected from the following:

and R^{27} is defined as above.

23. (original) The compound according to claim 22, wherein E is selected from -O-, -S(O)$_{0-2}$-, and -NH-; and G is -CH_2-.

24. (original) The compound according to claim 22, wherein E is either -CH_2- or -NH-; and G is selected from -O-, -S-, and -NH-.

25. (currently amended) The compound according ~~to either claim 23 or claim 24~~to claim 23, wherein each of R^3 is independently selected from -H, halogen, trihalomethyl, -OR^4, -CO_2R^4, -C(=O)R^4, and optionally substituted C_{1-6}alkyl.

26. (original) The compound according to claim 25, wherein at least one of R^{30} is a trifluoromethyl radical *meta*- to -E-G-C(=O)N(H)-.

27. (currently amended) The compound according ~~to either claim 1 or claim 16~~to claim 1, selected from Table 3:

Table 3

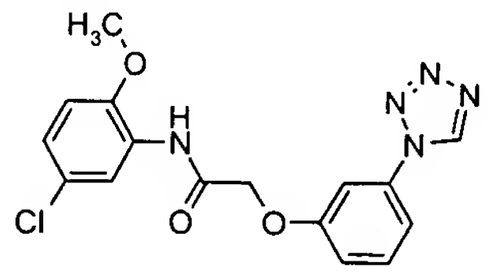
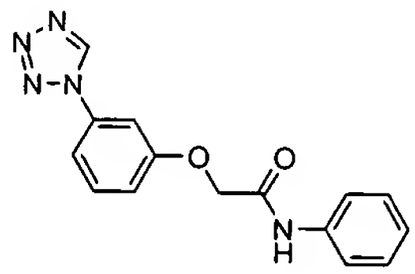
Entry	Name	Structure
1	N-[5-chloro-2-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
2	N-phenyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3

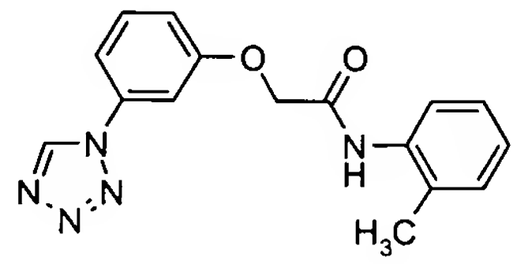
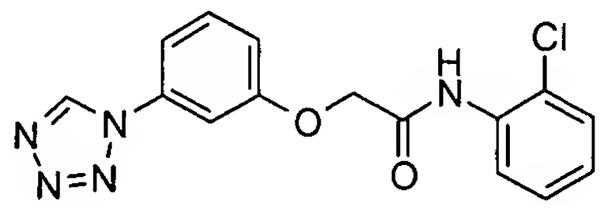
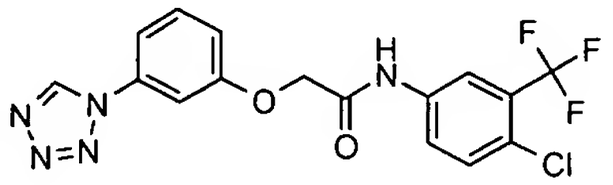
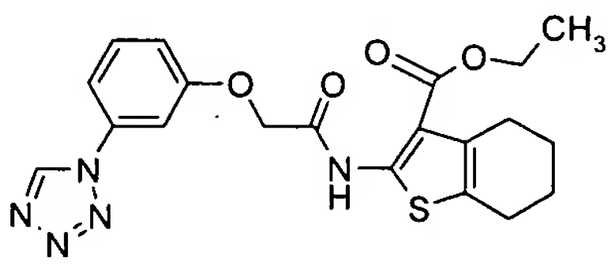
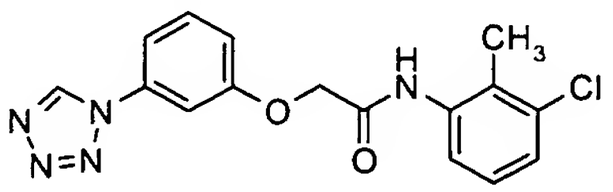
Entry	Name	Structure
3	N-(2-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
4	N-(2-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
5	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
6	ethyl 2-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]-4,5,6,7-tetrahydro-1-benzothiophene-3-carboxylate	
7	N-(3-chloro-2-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3

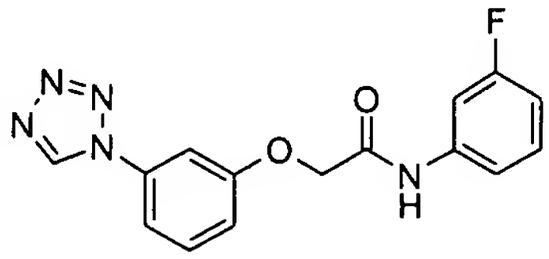
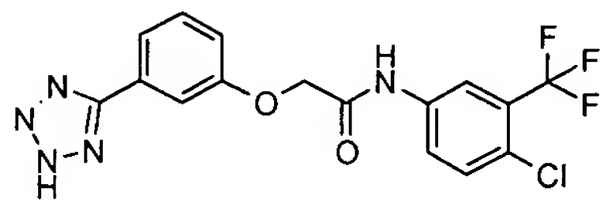
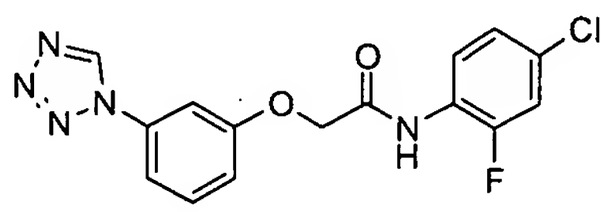
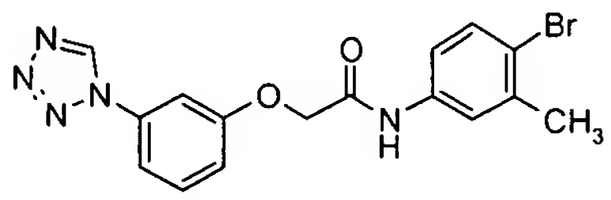
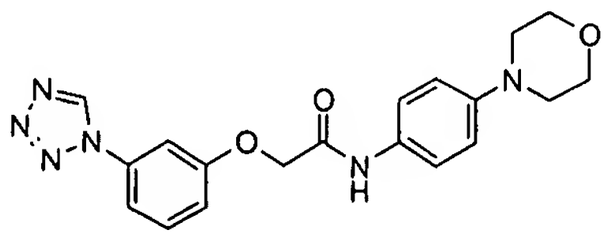
Entry	Name	Structure
8	N-(3-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
9	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(2H-tetrazol-5-yl)phenyl]oxy}acetamide	
10	N-(4-chloro-2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
11	N-(4-bromo-3-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
12	N-(4-morpholin-4-ylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3

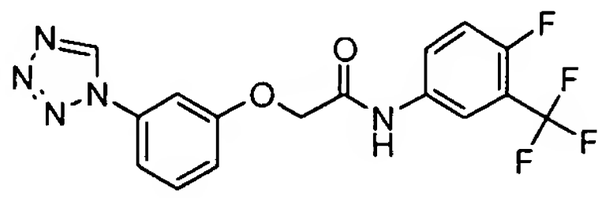
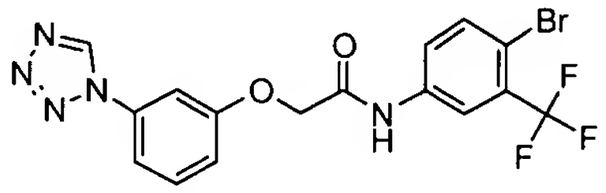
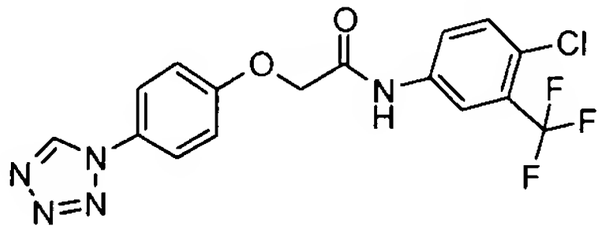
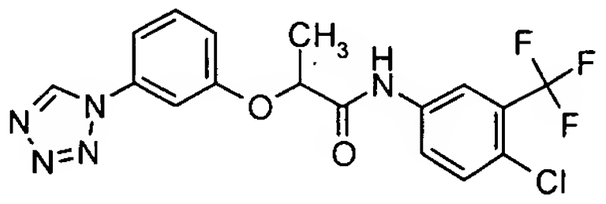
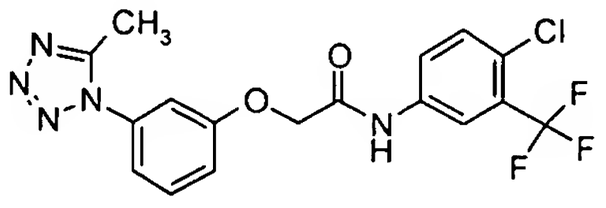
Entry	Name	Structure
13	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
14	N-[4-bromo-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
15	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
16	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	
17	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(5-methyl-1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3

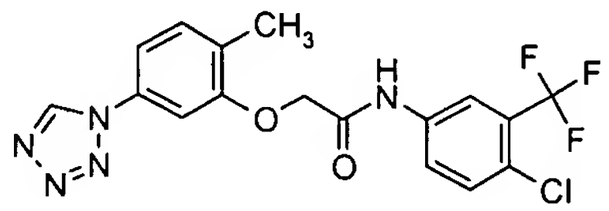
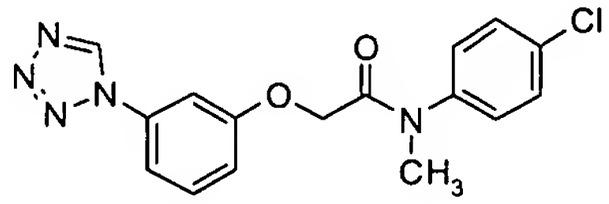
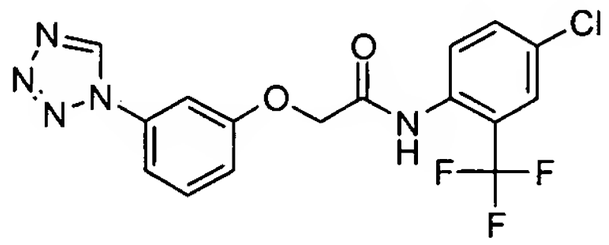
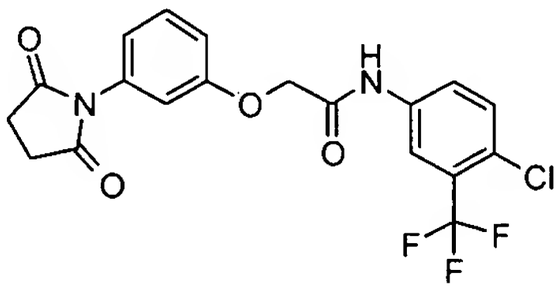
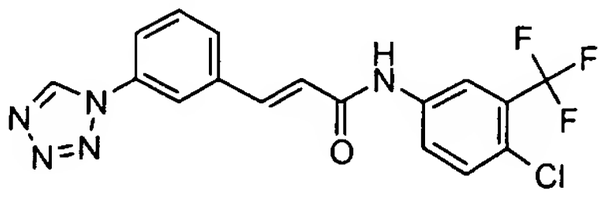
Entry	Name	Structure
18	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2-methyl-5-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
19	N-(4-chlorophenyl)-N-methyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
20	N-[4-chloro-2-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
21	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(2,5-dioxopyrrolidin-1-yl)phenyl]oxy}acetamide	
22	(2E)-N-[4-chloro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]prop-2-enamide	

Table 3

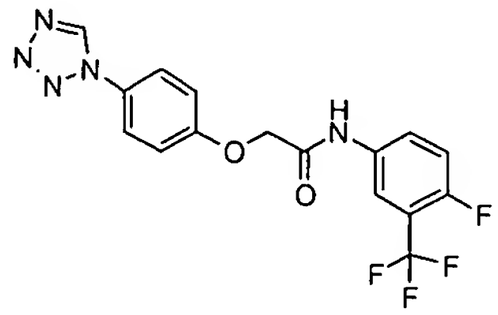
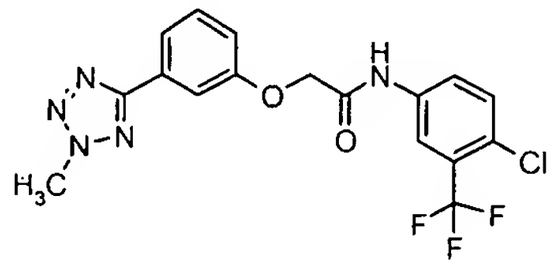
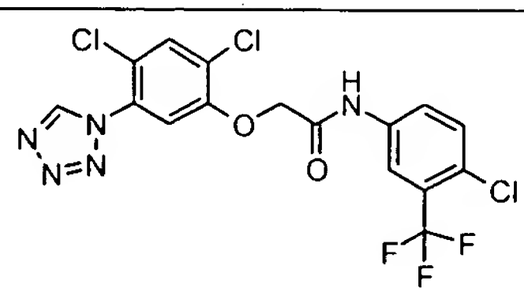
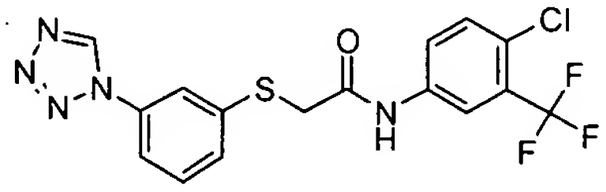
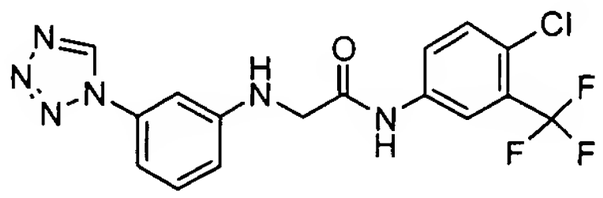
Entry	Name	Structure
23	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
24	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(2-methyl-2H-tetrazol-5-yl)phenyl]oxy}acetamide	
25	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2,4-dichloro-5-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
26	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]thio}acetamide	
27	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	

Table 3

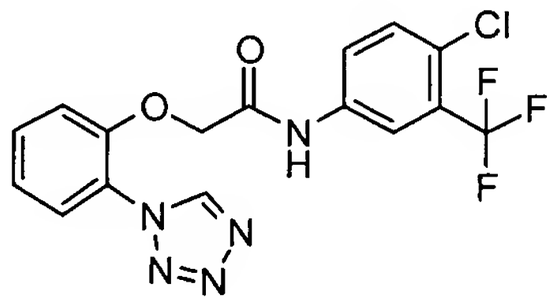
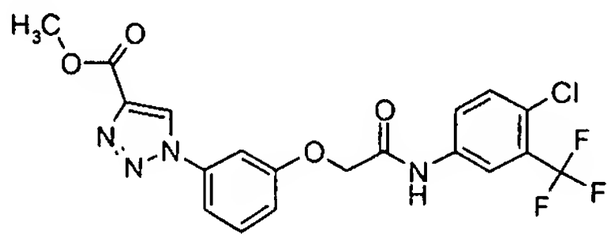
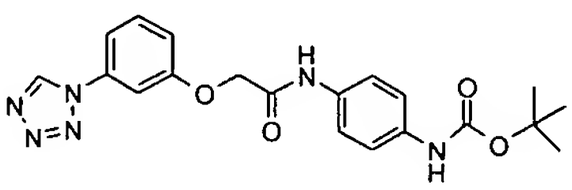
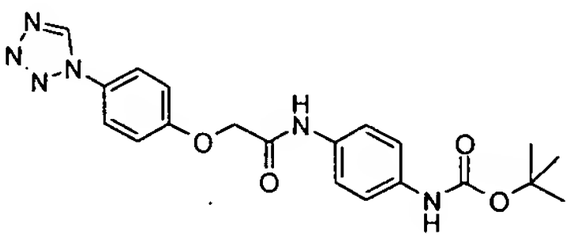
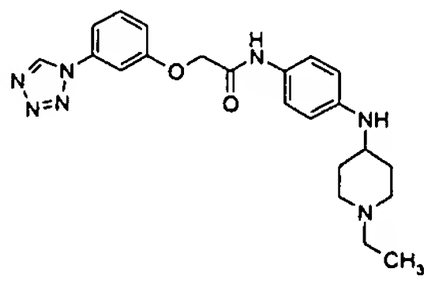
Entry	Name	Structure
28	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
29	methyl 1-{3-[(2-{[4-chloro-3-(trifluoromethyl)phenyl]amino}-2-oxoethyl)oxy]phenyl}-1H-1,2,3-triazole-4-carboxylate	
30	1,1-dimethylethyl {4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]phenyl}carbamate	
31	1,1-dimethylethyl {4-[({[4-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]phenyl}carbamate	
32	N-{4-[(1-ethylpiperidin-4-yl)amino]phenyl}-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3

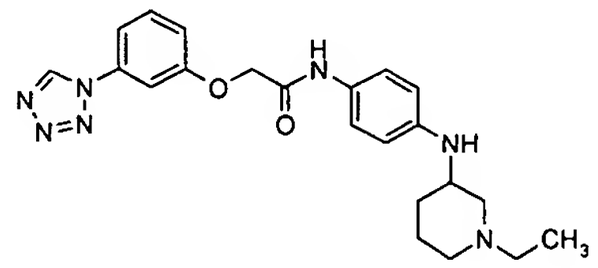
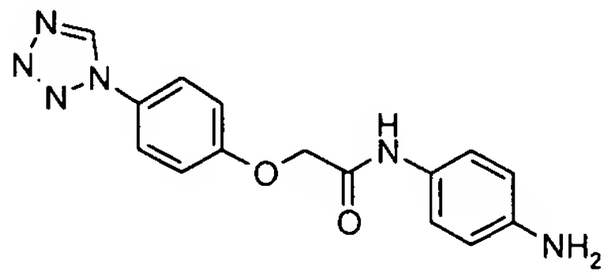
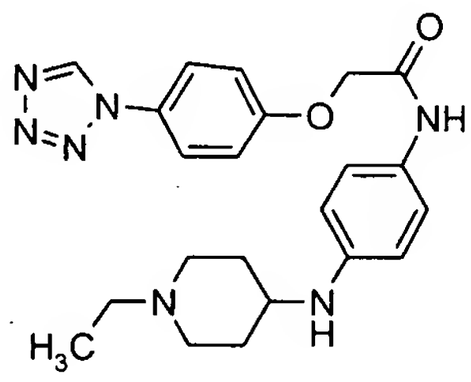
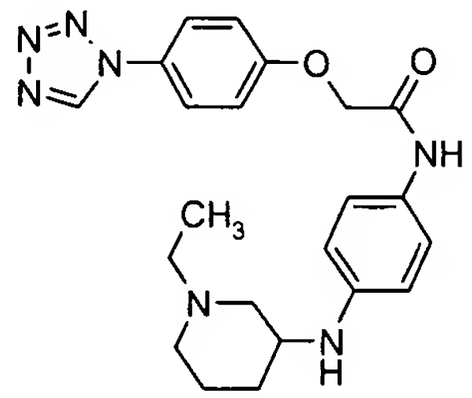
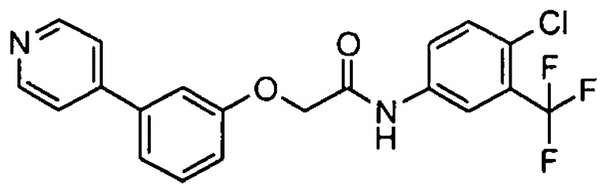
Entry	Name	Structure
33	N-{4-[(1-ethylpiperidin-3-yl)amino]phenyl}-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
34	N-(4-aminophenyl)-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
35	N-{4-[(1-ethylpiperidin-4-yl)amino]phenyl}-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
36	N-{4-[(1-ethylpiperidin-3-yl)amino]phenyl}-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
37	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyridin-4-ylphenyl)oxy]acetamide	

Table 3

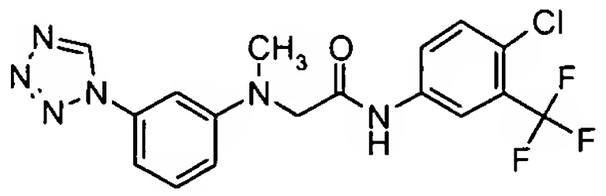
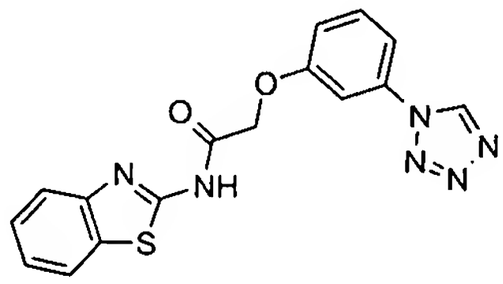
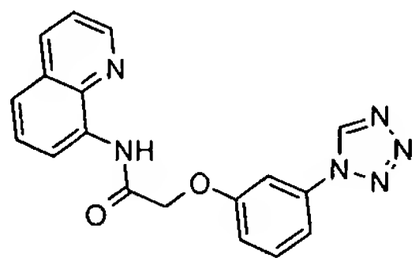
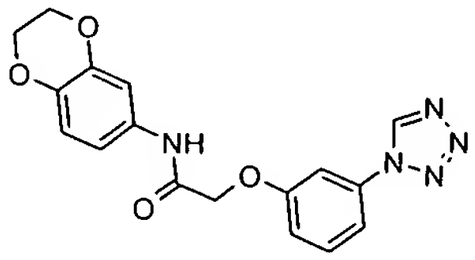
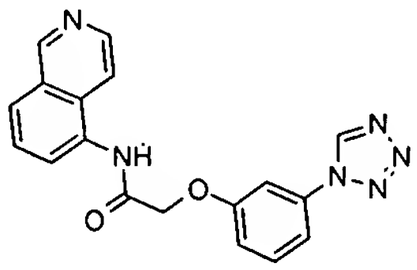
Entry	Name	Structure
38	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-methyl-N~2~-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
39	N-1,3-benzothiazol-2-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
40	N-quinolin-8-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
41	N-(2,3-dihydro-1,4-benzodioxin-6-yl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
42	N-isoquinolin-5-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3

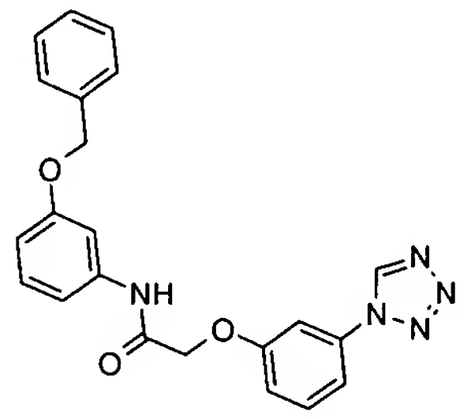
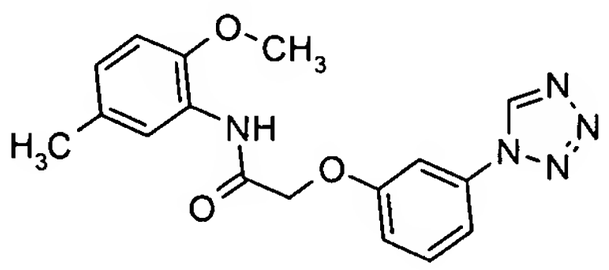
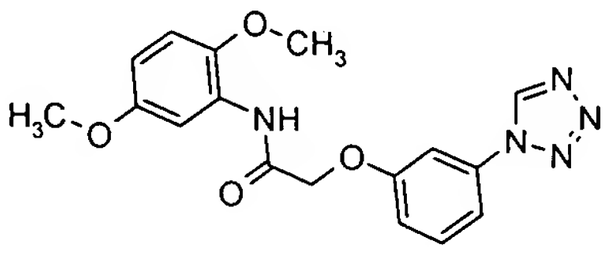
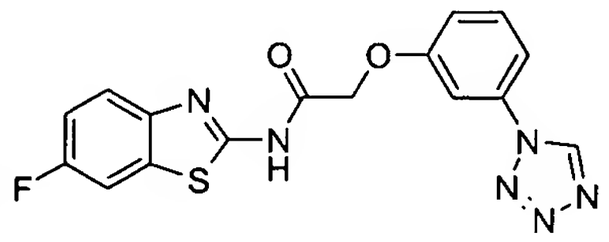
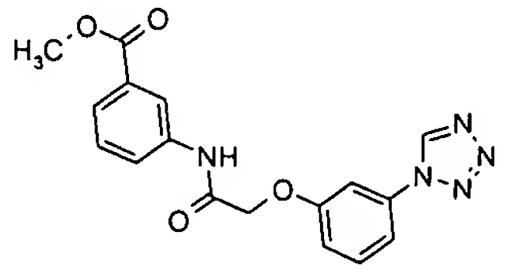
Entry	Name	Structure
43	N-{3-[(phenylmethyl)oxy]phenyl}-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
44	N-[5-methyl-2-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
45	N-[2,5-bis(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
46	N-(6-fluoro-1,3-benzothiazol-2-yl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
47	methyl 3-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoate	

Table 3

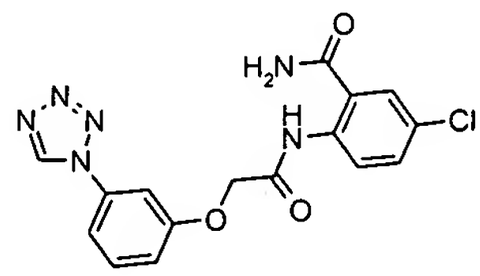
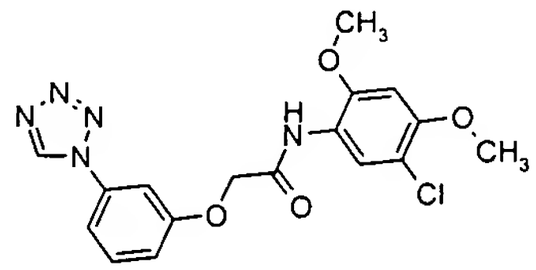
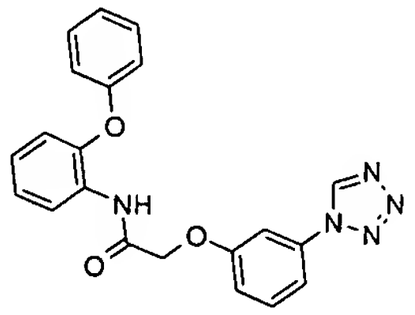
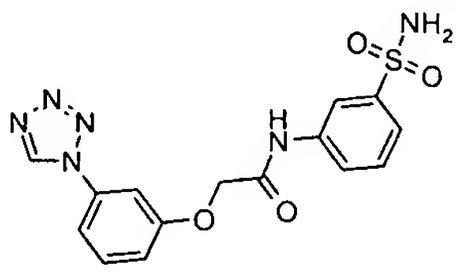
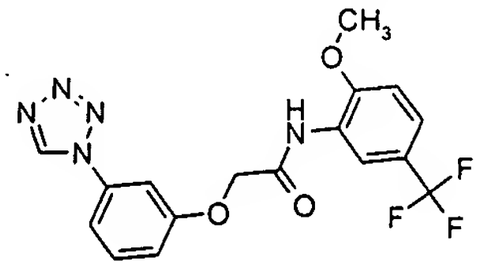
Entry	Name	Structure
48	5-chloro-2-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzamide	
49	N-[5-chloro-2,4-bis(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
50	N-[2-(phenyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
51	N-[3-(aminosulfonyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
52	N-[2-(methyloxy)-5-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3

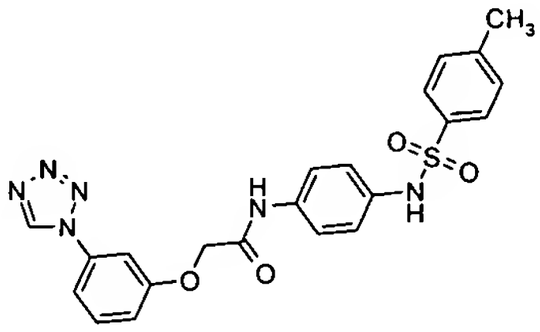
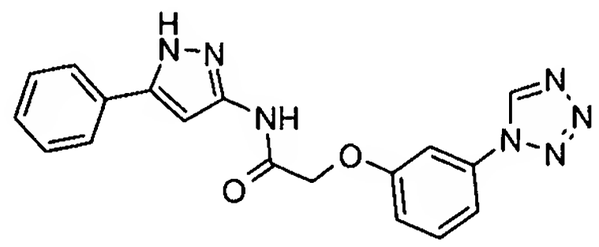
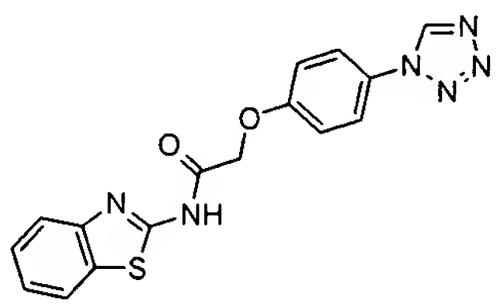
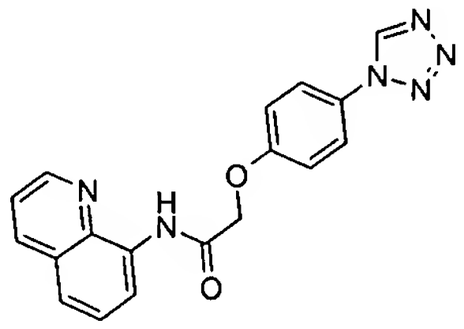
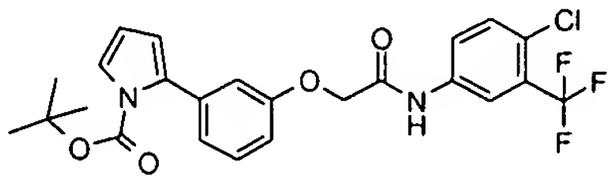
Entry	Name	Structure
53	N-(4-{[(4-methylphenyl)sulfonyl]amino}phenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
54	N-(5-phenyl-1H-pyrazol-3-yl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
55	N-1,3-benzothiazol-2-yl-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
56	N-quinolin-8-yl-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
57	1,1-dimethylethyl 2-{3-[(2-{[4-chloro-3-(trifluoromethyl)phenyl]amino}-2-oxoethyl)oxy]phenyl}-1H-pyrrole-1-carboxylate	

Table 3

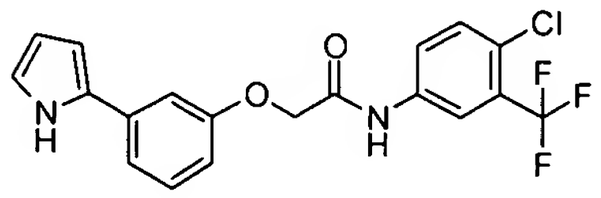
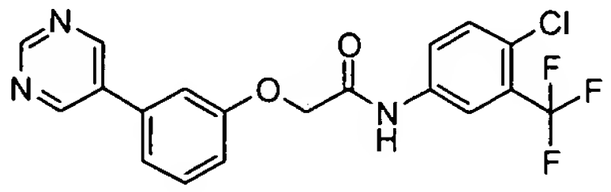
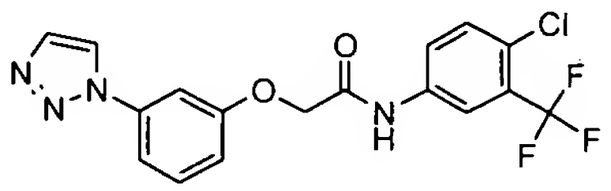
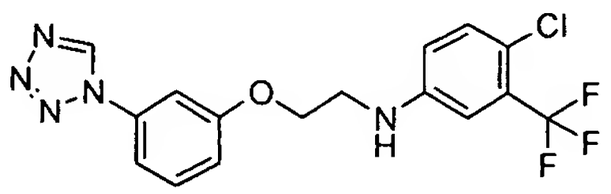
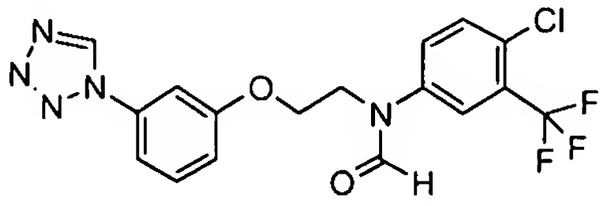
Entry	Name	Structure
58	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-pyrrol-2-yl)phenyl]oxy}acetamide	
59	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyrimidin-5-ylphenyl)oxy]acetamide	
60	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-1,2,3-triazol-1-yl)phenyl]oxy}acetamide	
61	4-chloro-N-(2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}ethyl)-3-(trifluoromethyl)aniline	
62	N-[4-chloro-3-(trifluoromethyl)phenyl]-N-(2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}ethyl)formamide	

Table 3

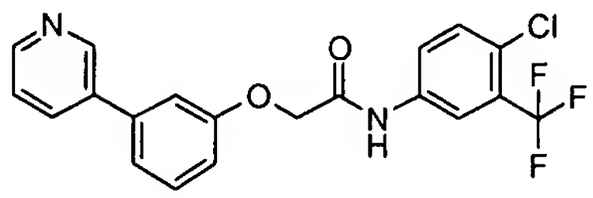
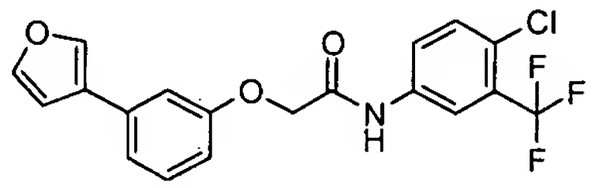
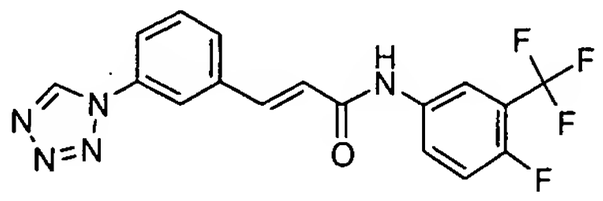
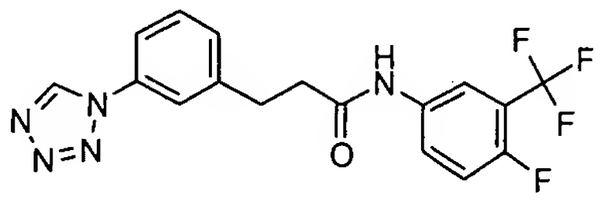
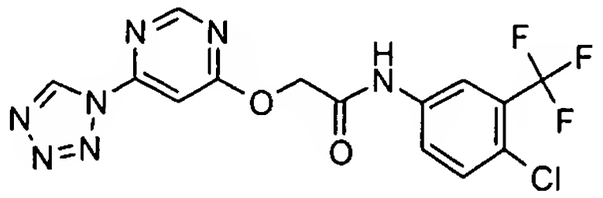
Entry	Name	Structure
63	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyridin-3-ylphenyl)oxy]acetamide	
64	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-furan-3-ylphenyl)oxy]acetamide	
65	(2E)-N-[4-fluoro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]prop-2-enamide	
66	N-[4-fluoro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]propanamide	
67	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[6-(1H-tetrazol-1-yl)pyrimidin-4-yl]oxy}acetamide	

Table 3

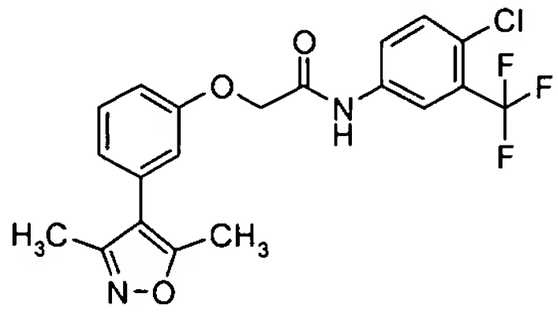
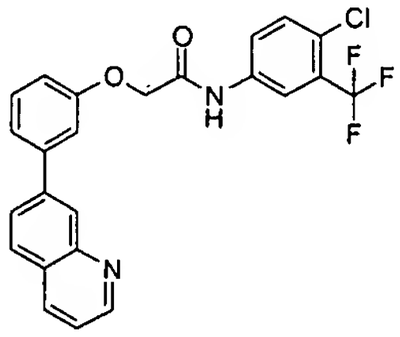
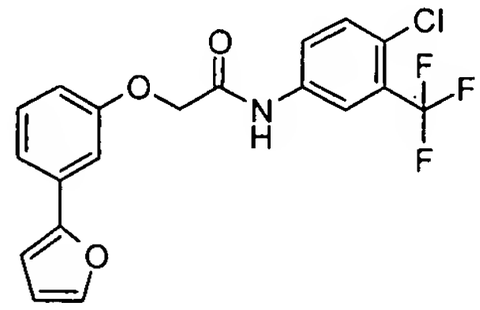
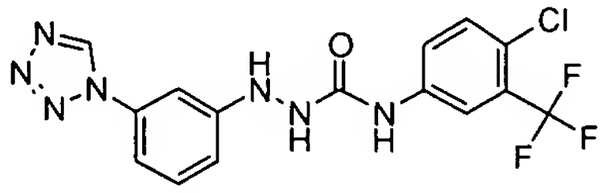
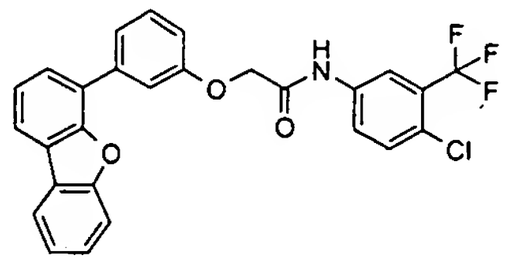
Entry	Name	Structure
68	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(3,5-dimethylisoxazol-4-yl)phenyl]oxy}acetamide	
69	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-quinolin-7-ylphenyl)oxy]acetamide	
70	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-furan-2-ylphenyl)oxy]acetamide	
71	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
72	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-dibenzo[b,d]furan-4-ylphenyl)oxy]acetamide	

Table 3

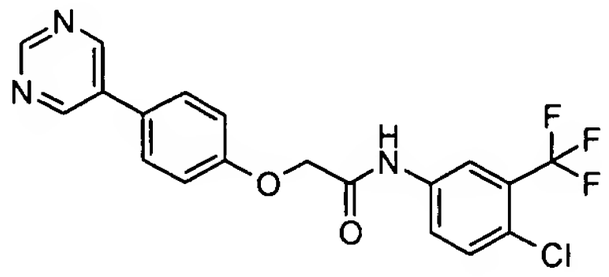
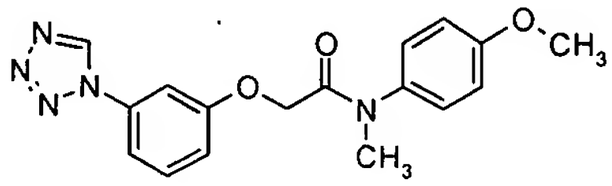
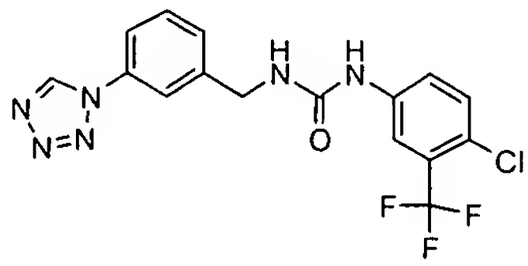
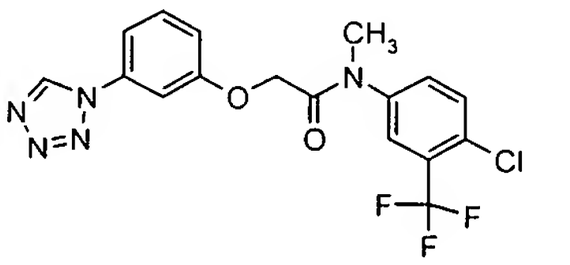
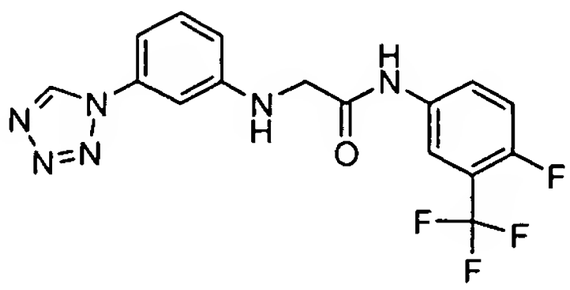
Entry	Name	Structure
73	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	
74	N-methyl-N-[4-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
75	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(1H-tetrazol-1-yl)phenyl]methyl}urea	
76	N-[4-chloro-3-(trifluoromethyl)phenyl]-N-methyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
77	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N~2~-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	

Table 3

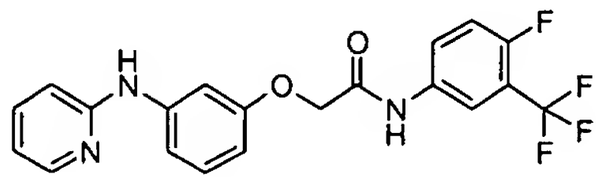
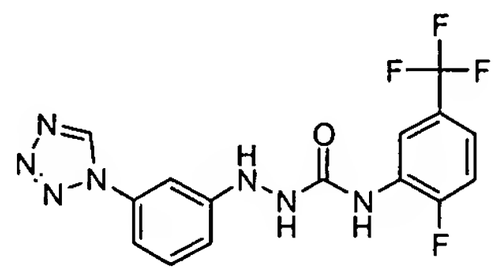
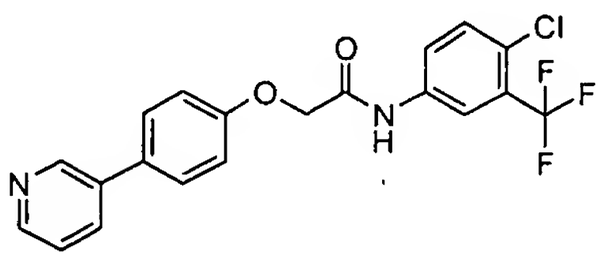
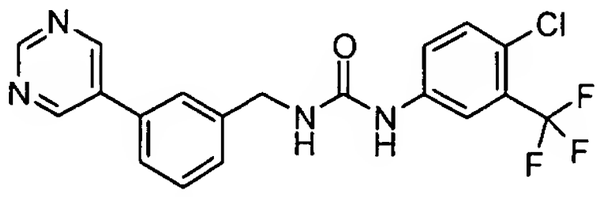
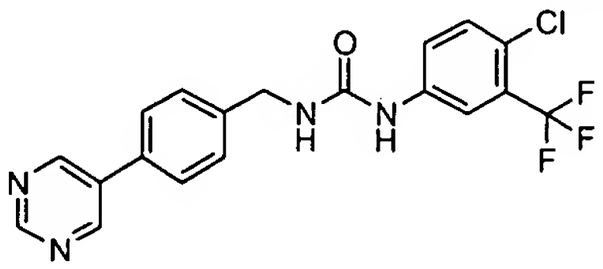
Entry	Name	Structure
78	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-{[3-(pyridin-2-ylamino)phenyl]oxy}acetamide	
79	N-[2-fluoro-5-(trifluoromethyl)phenyl]-2-[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
80	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide	
81	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyrimidin-5-ylphenyl)methyl]urea	
82	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyrimidin-5-ylphenyl)methyl]urea	

Table 3

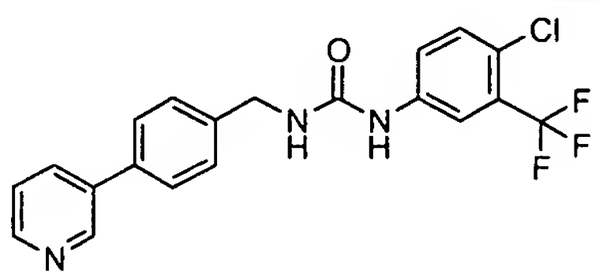
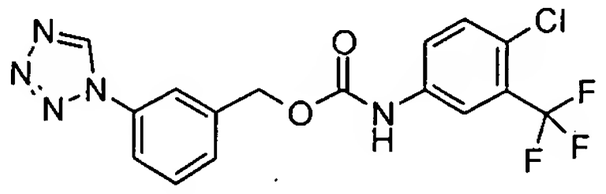
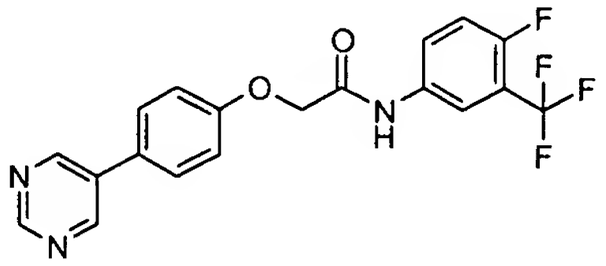
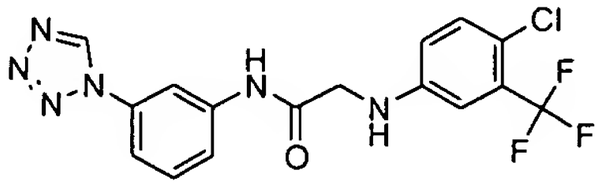
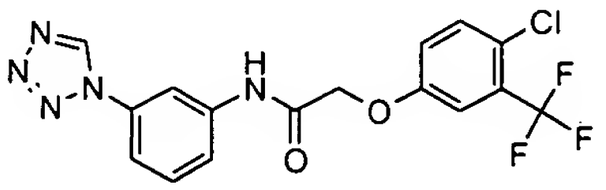
Entry	Name	Structure
83	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	
84	[3-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
85	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	
86	N~2~-[4-chloro-3-(trifluoromethyl)phenyl]-N-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
87	2-{[4-chloro-3-(trifluoromethyl)phenyl]oxy}-N-[3-(1H-tetrazol-1-yl)phenyl]acetamide	

Table 3

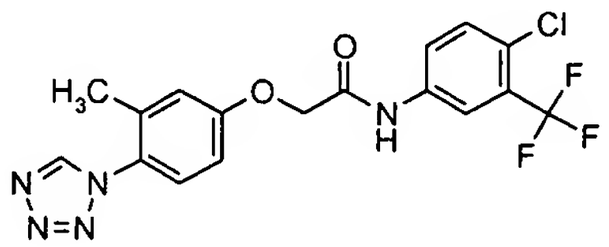
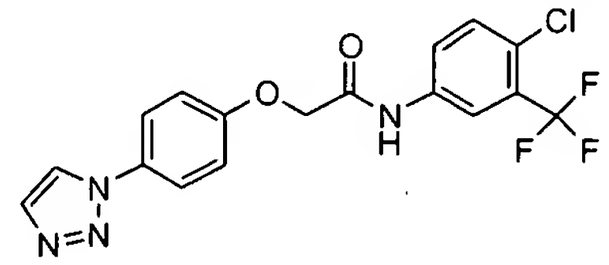
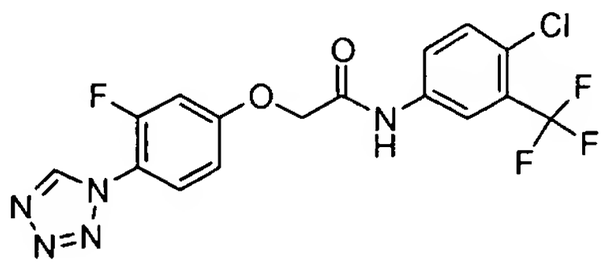
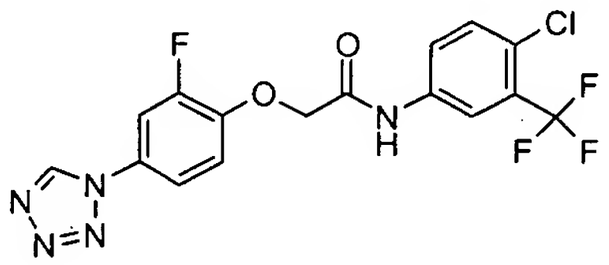
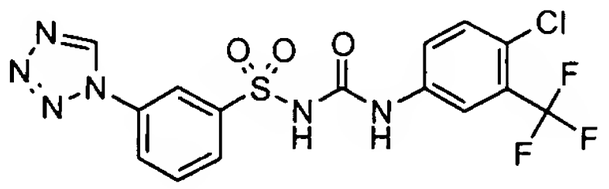
Entry	Name	Structure
88	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-methyl-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
89	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-1,2,3-triazol-1-yl)phenyl]oxy}acetamide	
90	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-fluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
91	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2-fluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
92	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-3-(1H-tetrazol-1-yl)benzenesulfonamide	

Table 3

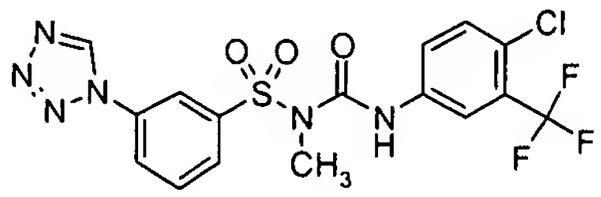
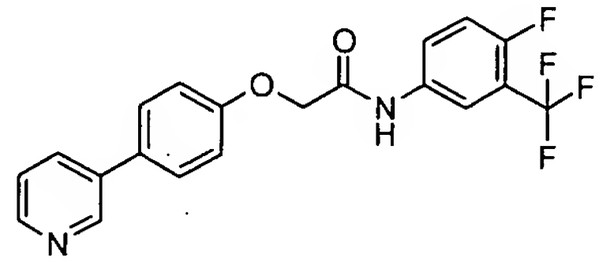
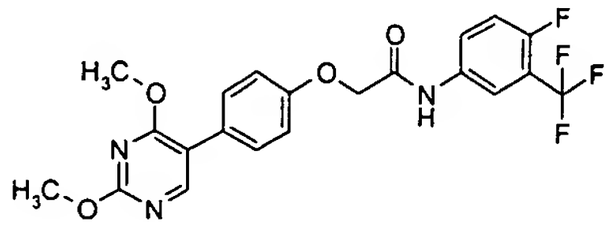
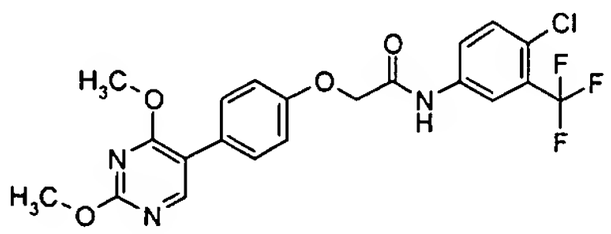
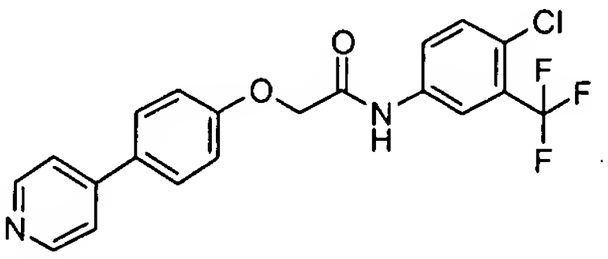
Entry	Name	Structure
93	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-N-methyl-3-(1H-tetrazol-1-yl)benzenesulfonamide	
94	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide	
95	2-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}oxy)-N-[4-fluoro-3-(trifluoromethyl)phenyl]acetamide	
96	2-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}oxy)-N-[4-chloro-3-(trifluoromethyl)phenyl]acetamide	
97	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-4-ylphenyl)oxy]acetamide	

Table 3

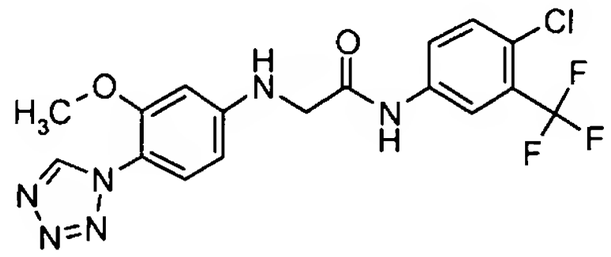
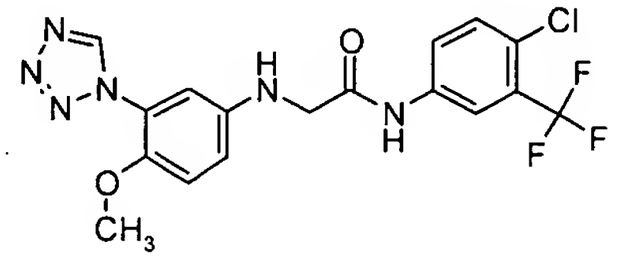
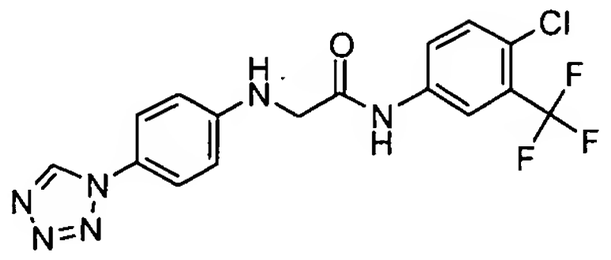
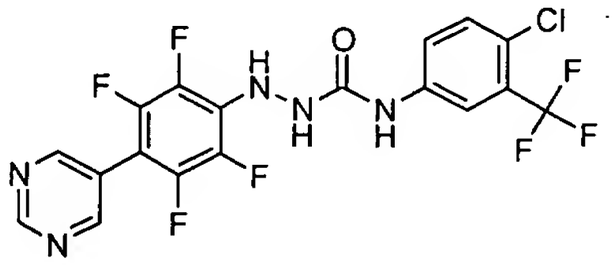
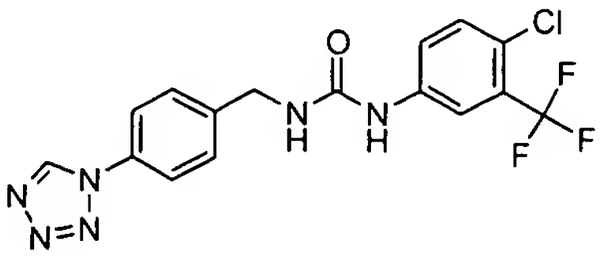
Entry	Name	Structure
98	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(methyloxy)-4-(1H-tetrazol-1-yl)phenyl]glycinamide	
99	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[4-(methyloxy)-3-(1H-tetrazol-1-yl)phenyl]glycinamide	
100	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[4-(1H-tetrazol-1-yl)phenyl]glycinamide	
101	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(2,3,5,6-tetrafluoro-4-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
102	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(1H-tetrazol-1-yl)phenyl]methyl}urea	

Table 3

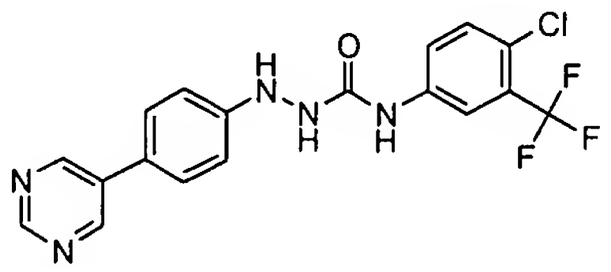
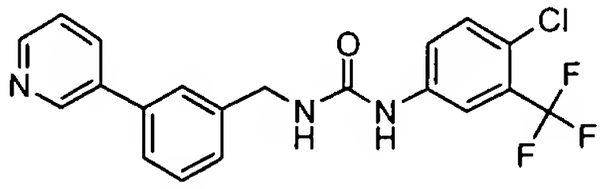
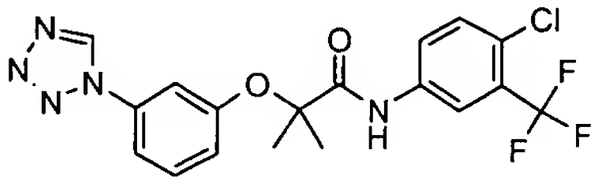
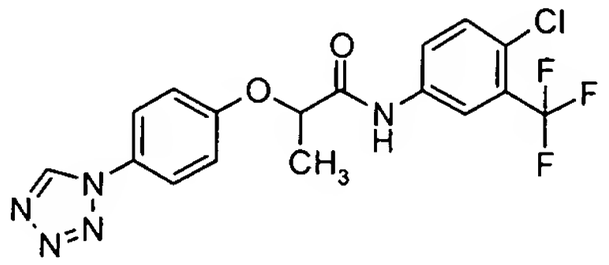
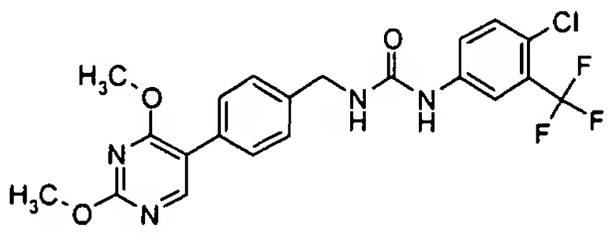
Entry	Name	Structure
103	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
104	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyridin-3-ylphenyl)methyl]urea	
105	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-methyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	
106	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	
107	N-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Table 3

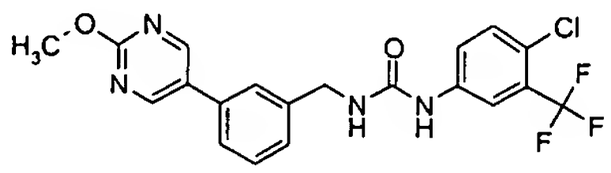
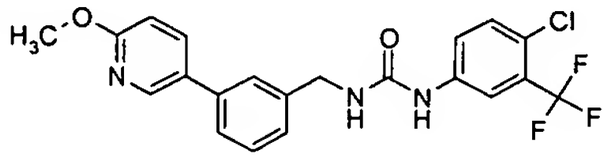
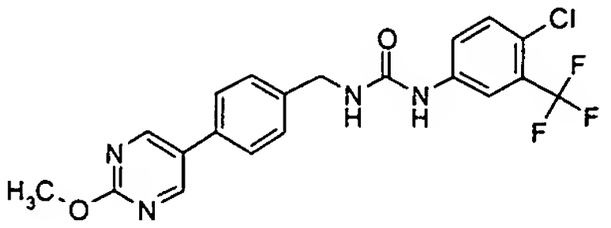
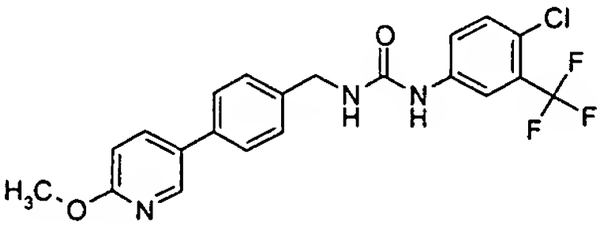
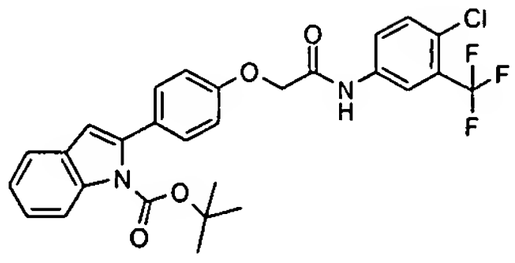
Entry	Name	Structure
108	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[2-(methyloxy)pyrimidin-5-yl]phenyl}methyl)urea	
109	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[6-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	
110	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[2-(methyloxy)pyrimidin-5-yl]phenyl}methyl)urea	
111	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[6-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	
112	1,1-dimethylethyl 2-{4-[(2-{[4-chloro-3-(trifluoromethyl)phenyl]amino}-2-oxoethyl)oxy]phenyl}-1H-indole-1-carboxylate	

Table 3

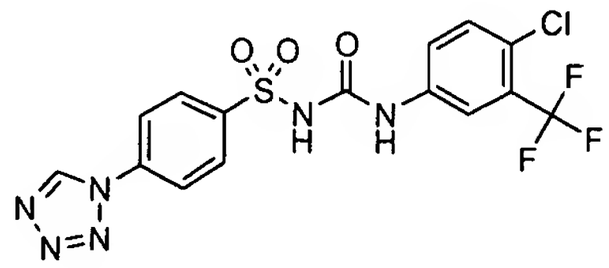
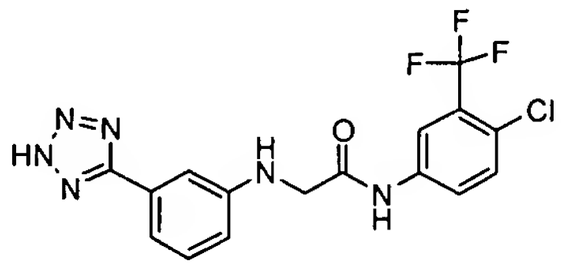
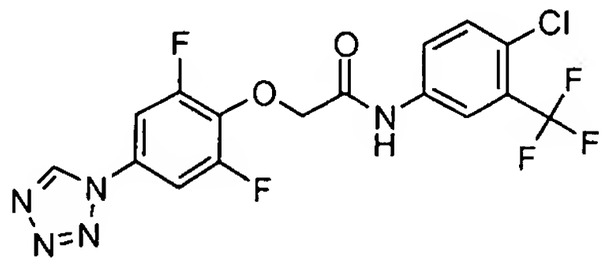
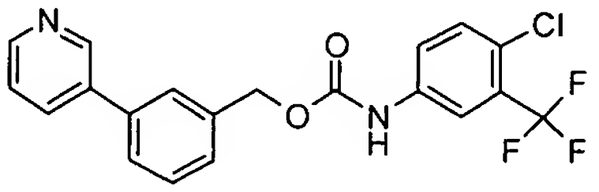
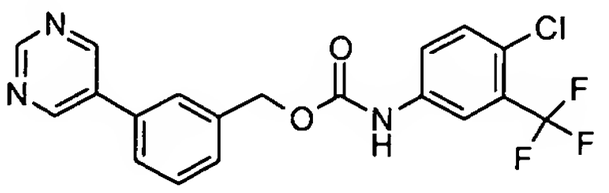
Entry	Name	Structure
113	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-4-(1H-tetrazol-1-yl)benzenesulfonamide	
114	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(2H-tetrazol-5-yl)phenyl]glycinamide	
115	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2,6-difluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
116	(3-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
117	(3-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

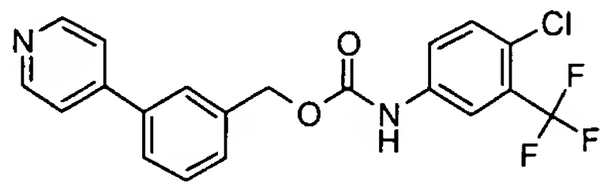
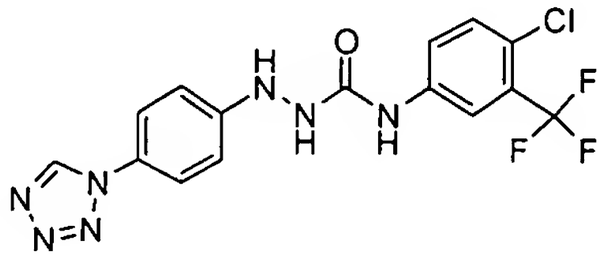
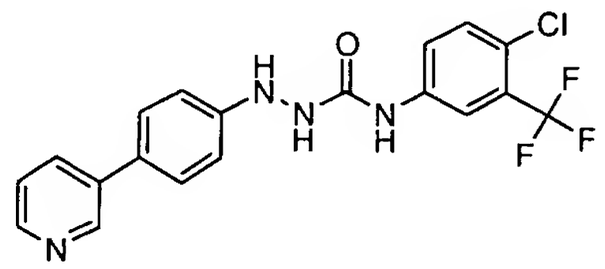
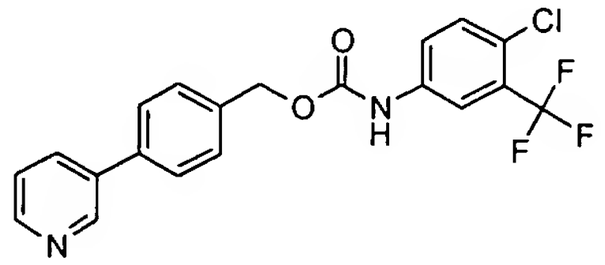
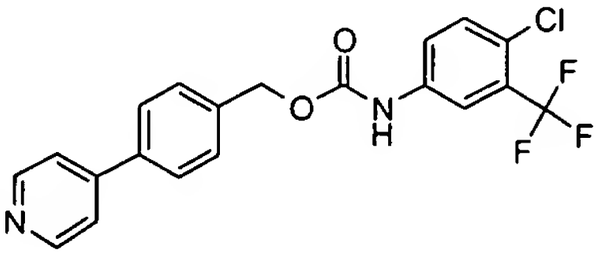
Entry	Name	Structure
118	(3-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
119	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[4-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
120	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-3-ylphenyl)hydrazinecarboxamide	
121	(4-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
122	(4-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

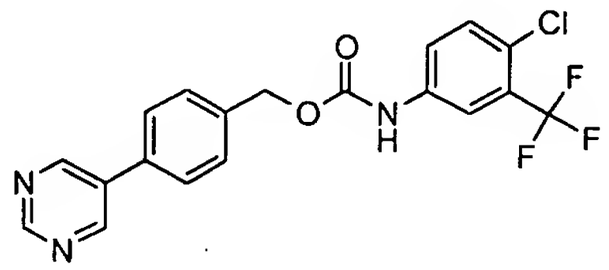
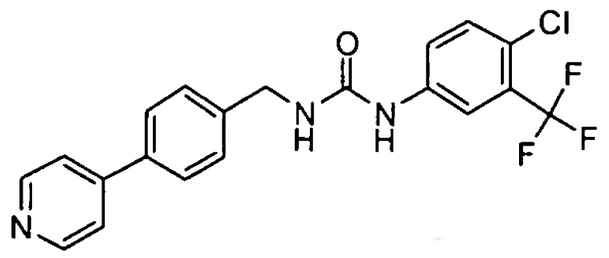
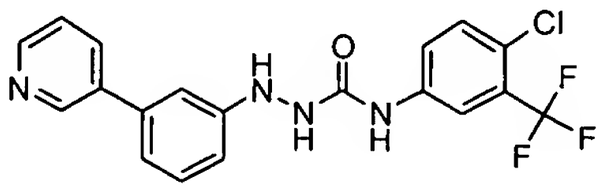
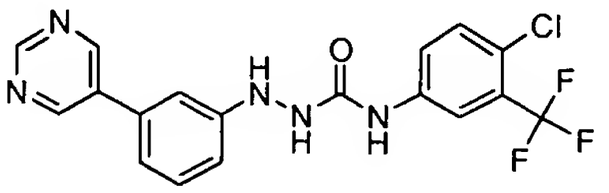
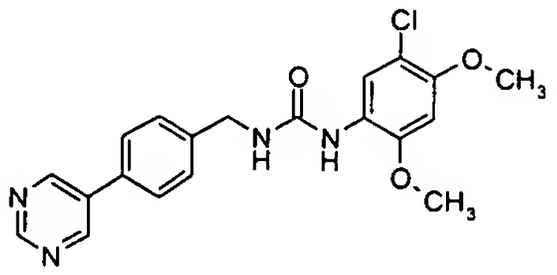
Entry	Name	Structure
123	(4-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
124	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-4-ylphenyl)methyl]urea	
125	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-3-ylphenyl)hydrazinecarboxamide	
126	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
127	N-[5-chloro-2,4-bis(methyloxy)phenyl]-N'-[(4-pyrimidin-5-ylphenyl)methyl]urea	

Table 3

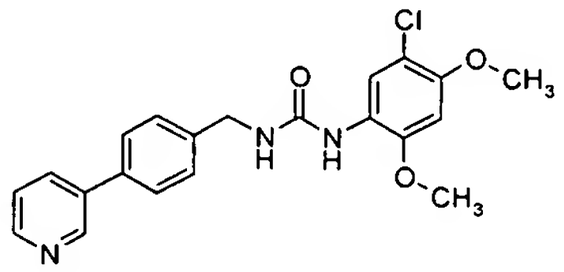
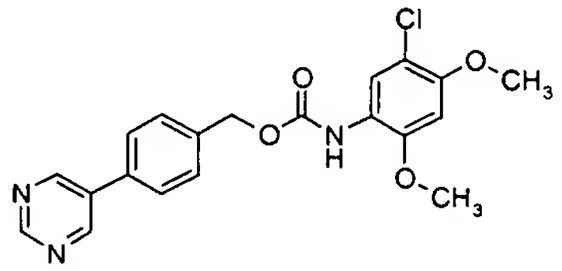
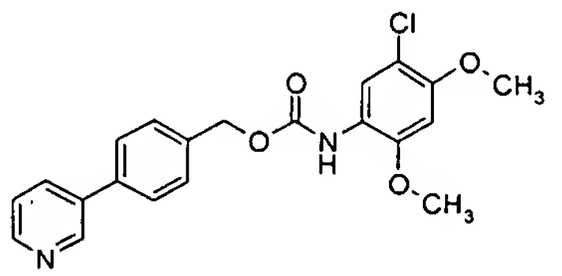
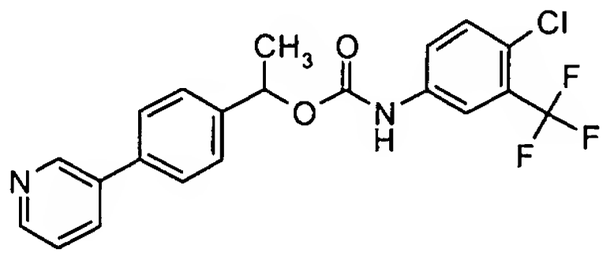
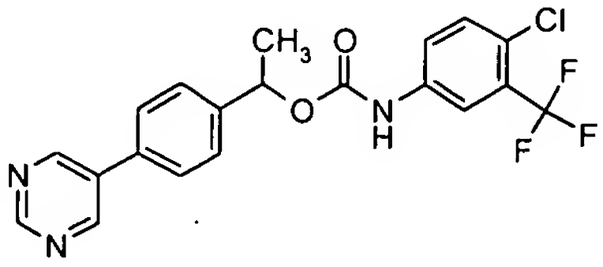
Entry	Name	Structure
128	N-[5-chloro-2,4-bis(methyloxy)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	
129	(4-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	
130	(4-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	
131	1-(4-pyridin-3-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
132	1-(4-pyrimidin-5-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

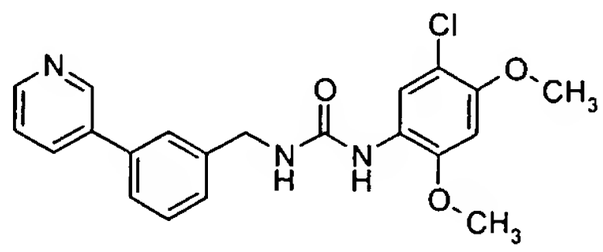
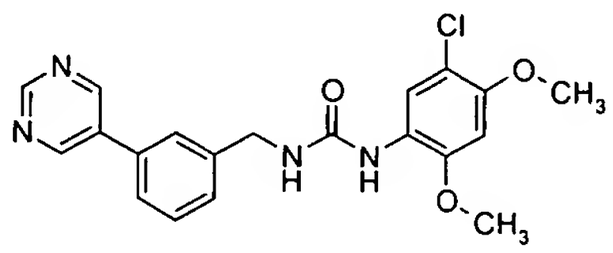
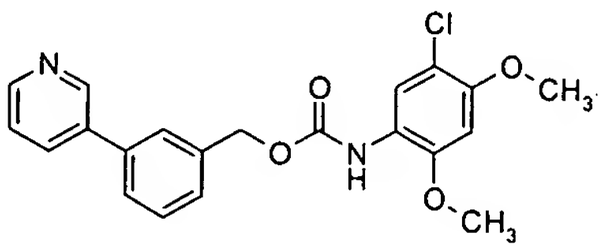
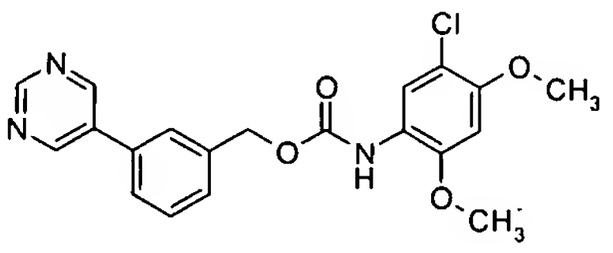
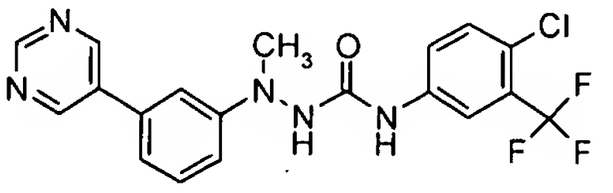
Entry	Name	Structure
133	N-[5-chloro-2,4-bis(methyloxy)phenyl]-N'-[(3-pyridin-3-ylphenyl)methyl]urea	
134	N-[5-chloro-2,4-bis(methyloxy)phenyl]-N'-[(3-pyrimidin-5-ylphenyl)methyl]urea	
135	(3-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	
136	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	
137	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-methyl-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide	

Table 3

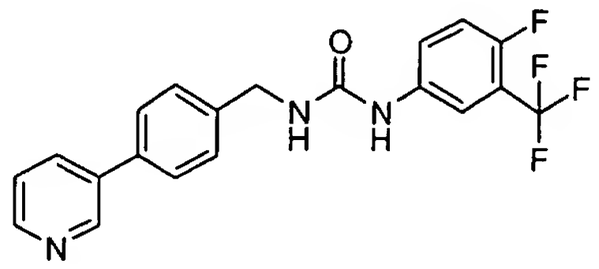
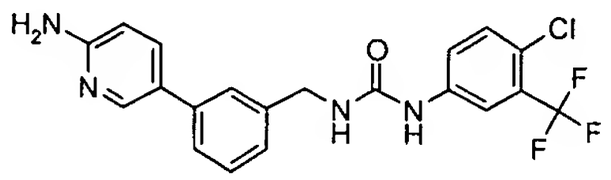
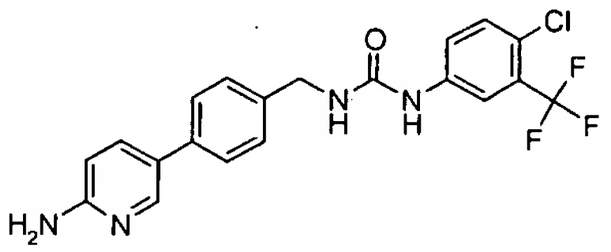
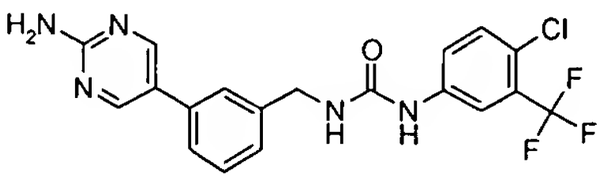
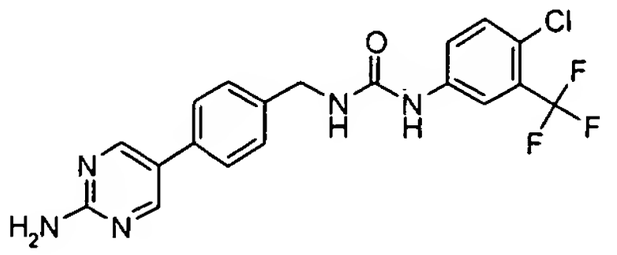
Entry	Name	Structure
138	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	
139	N-{[3-(6-aminopyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
140	N-{[4-(6-aminopyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
141	N-{[3-(2-aminopyrimidin-5-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
142	N-{[4-(2-aminopyrimidin-5-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Table 3

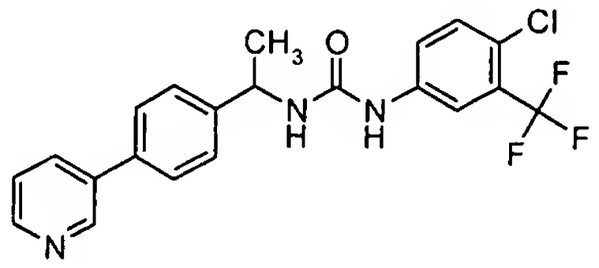
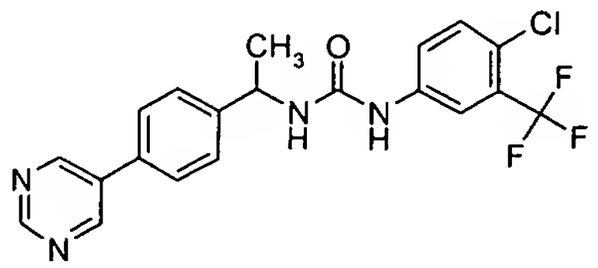
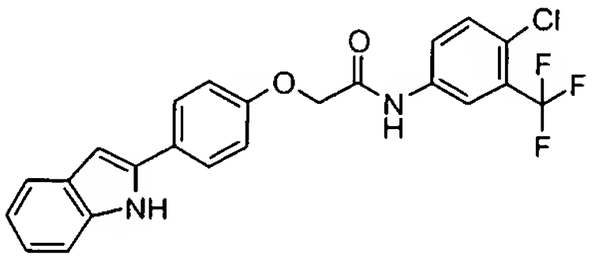
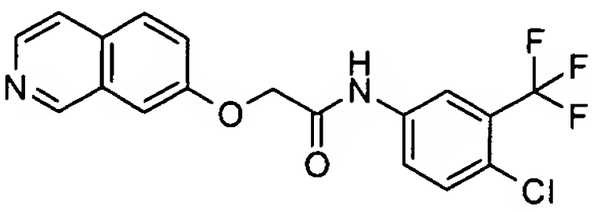
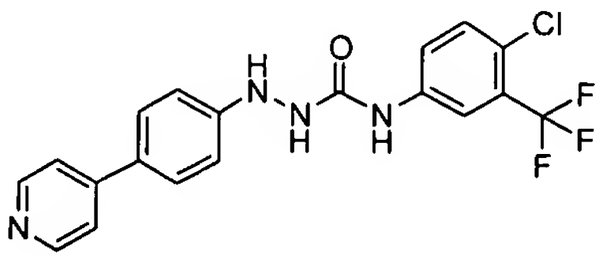
Entry	Name	Structure
143	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyridin-3-ylphenyl)ethyl]urea	
144	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyrimidin-5-ylphenyl)ethyl]urea	
145	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-indol-2-yl)phenyl]oxy}acetamide	
146	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(isoquinolin-7-yloxy)acetamide	
147	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-4-ylphenyl)hydrazinecarboxamide	

Table 3

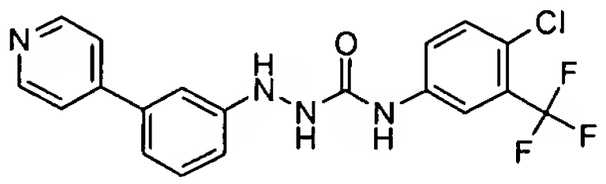
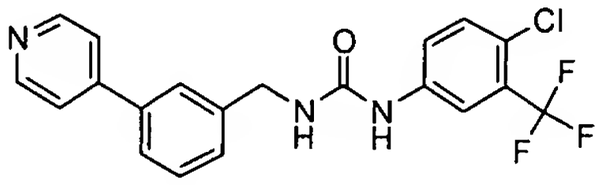
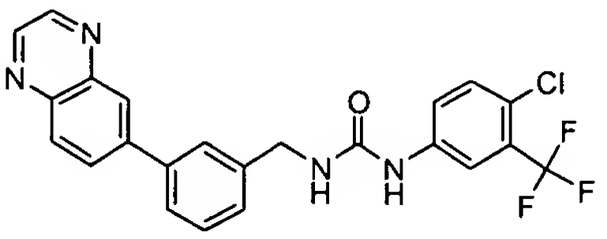
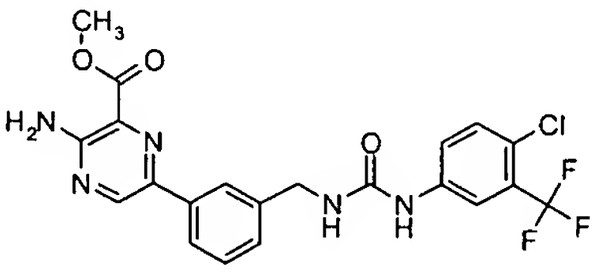
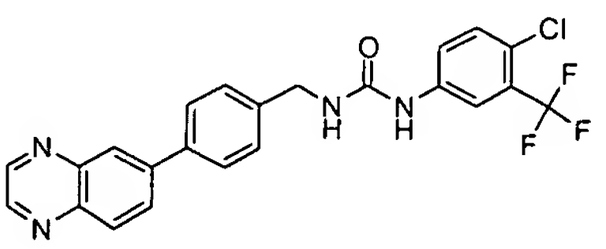
Entry	Name	Structure
148	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-4-ylphenyl)hydrazinecarboxamide	
149	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyridin-4-ylphenyl)methyl]urea	
150	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-quinoxalin-6-ylphenyl)methyl]urea	
151	methyl 3-amino-6-(3-{[({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)pyrazine-2-carboxylate	
152	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-quinoxalin-6-ylphenyl)methyl]urea	

Table 3

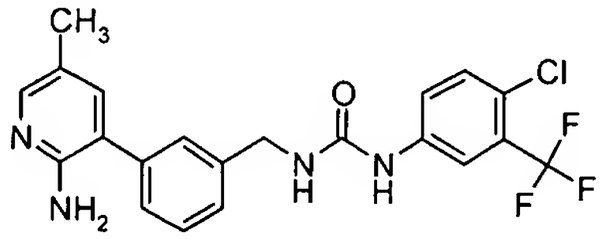
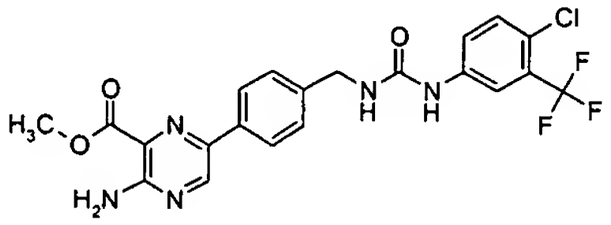
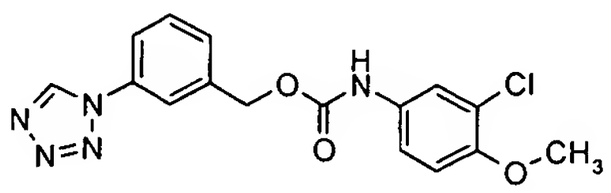
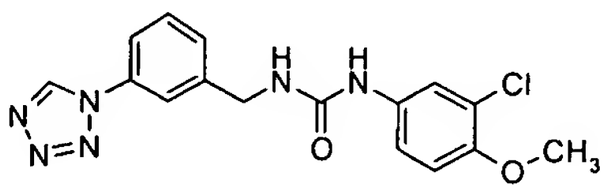
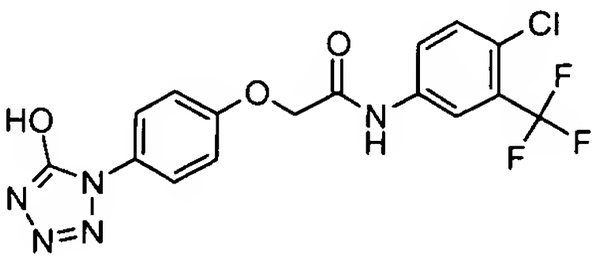
Entry	Name	Structure
153	N-{[3-(2-amino-5-methylpyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
154	methyl 3-amino-6-(4-{[({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)pyrazine-2-carboxylate	
155	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-chloro-4-(methyloxy)phenyl]carbamate	
156	N-[3-chloro-4-(methyloxy)phenyl]-N'-{[3-(1H-tetrazol-1-yl)phenyl]methyl}urea	
157	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(5-hydroxy-1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3

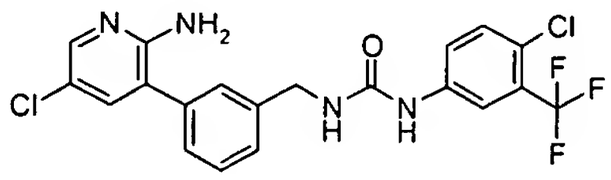
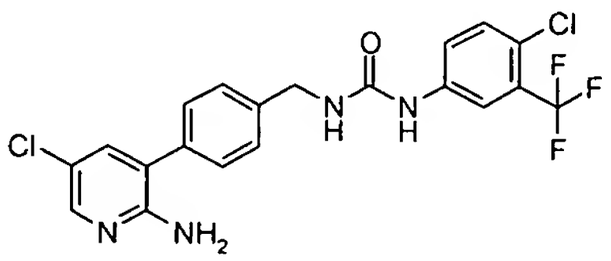
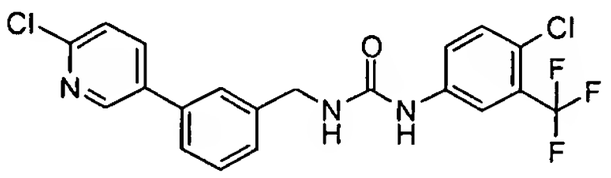
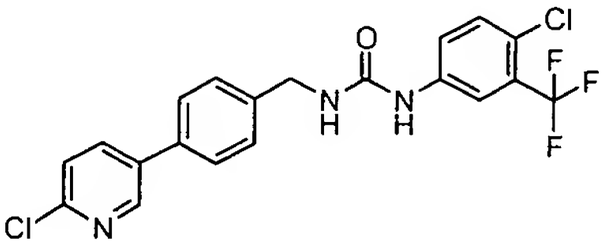
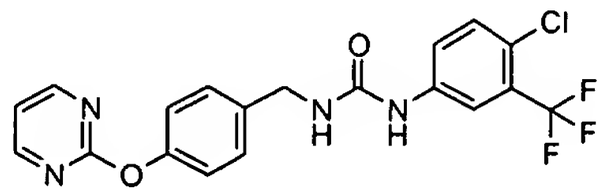
Entry	Name	Structure
158	N-{[3-(2-amino-5-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
159	N-{[4-(2-amino-5-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
160	N-{[3-(6-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
161	N-{[4-(6-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
162	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(pyrimidin-2-yloxy)phenyl]methyl}urea	

Table 3

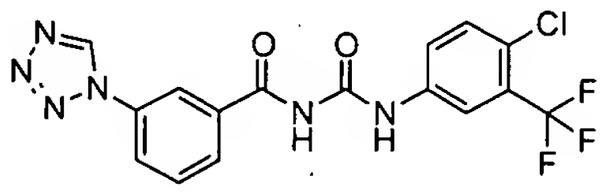
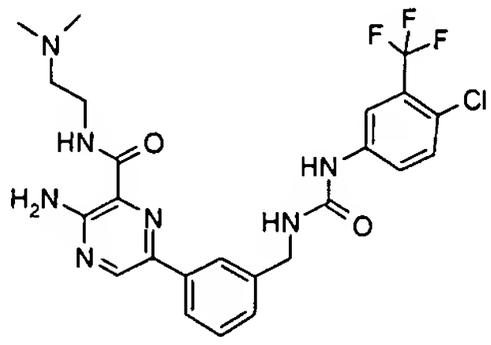
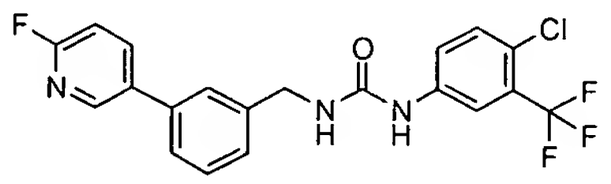
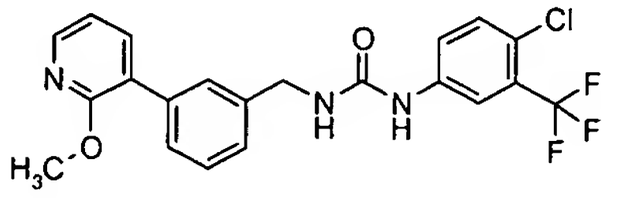
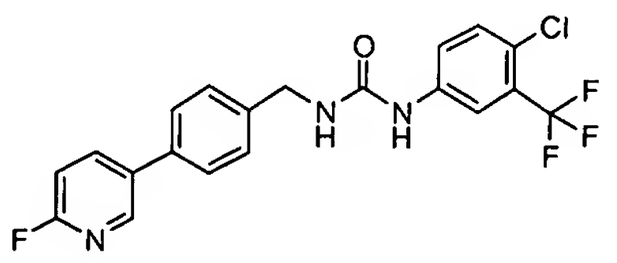
Entry	Name	Structure
163	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-3-(1H-tetrazol-1-yl)benzamide	
164	3-amino-6-(3-{[({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)-N-[2-(dimethylamino)ethyl]pyrazine-2-carboxamide	
165	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(6-fluoropyridin-3-yl)phenyl]methyl}urea	
166	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[2-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	
167	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(6-fluoropyridin-3-yl)phenyl]methyl}urea	

Table 3

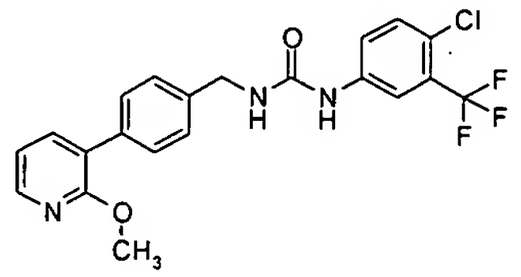
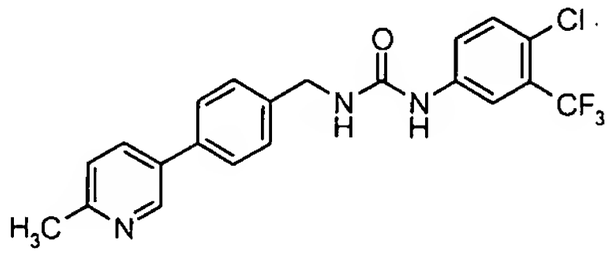
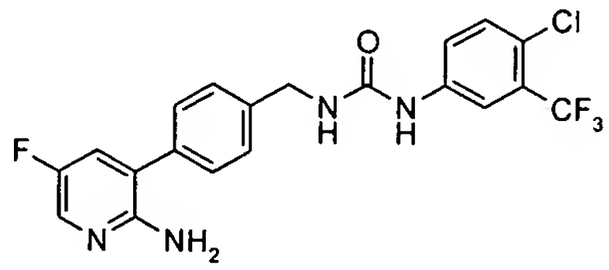
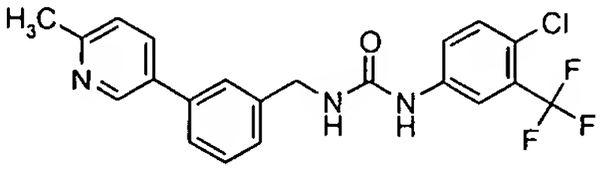
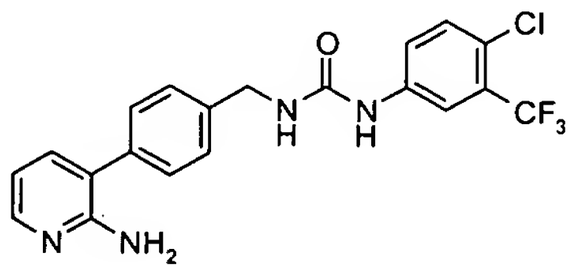
Entry	Name	Structure
168	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[2-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	
169	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(6-methylpyridin-3-yl)phenyl]methyl}urea	
170	N-{[4-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
171	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(6-methylpyridin-3-yl)phenyl]methyl}urea	
172	N-{[4-(2-aminopyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Table 3

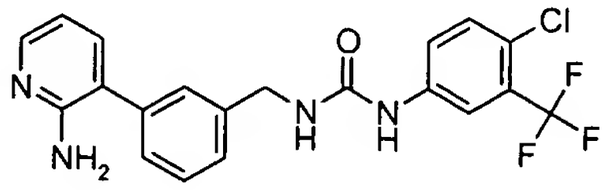
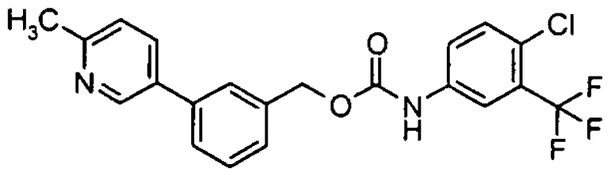
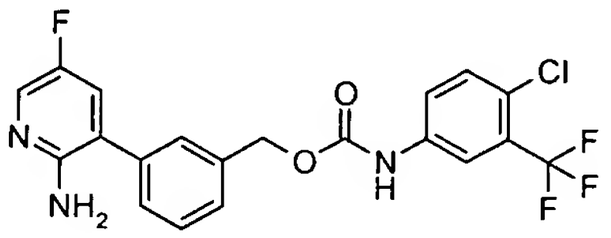
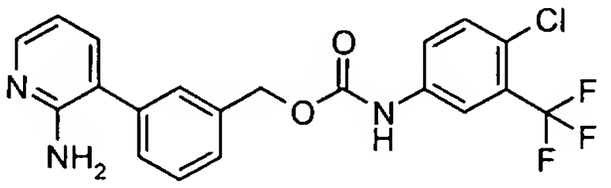
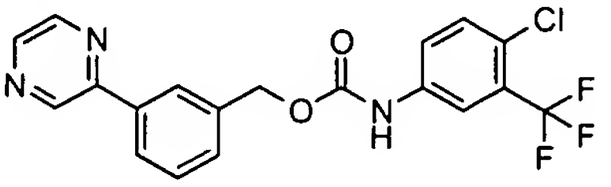
Entry	Name	Structure
173	N-{[3-(2-aminopyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
174	[3-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
175	[3-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
176	[3-(2-aminopyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
177	(3-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

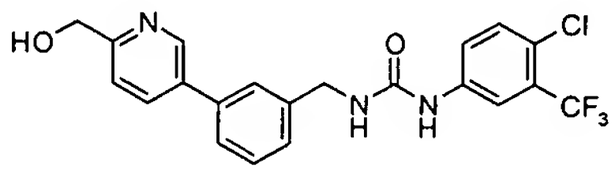
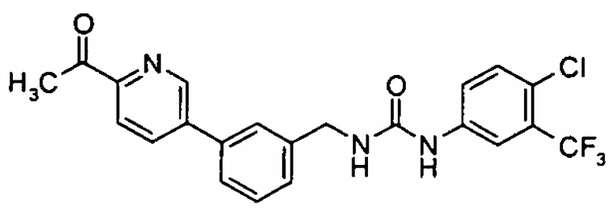
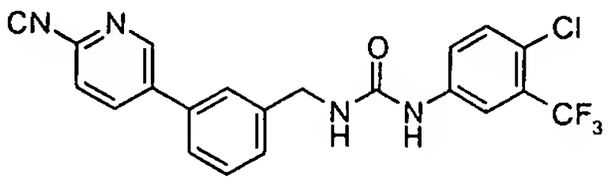
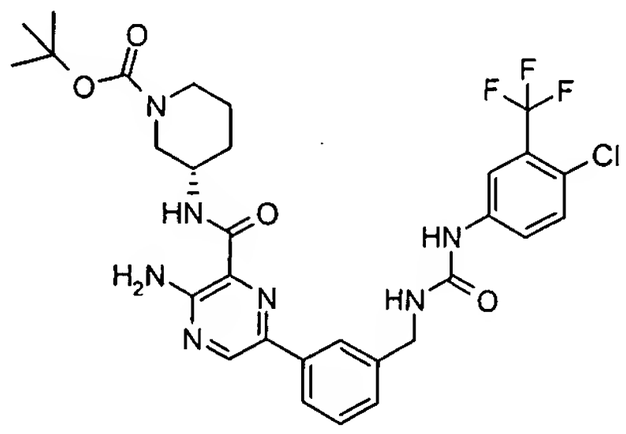
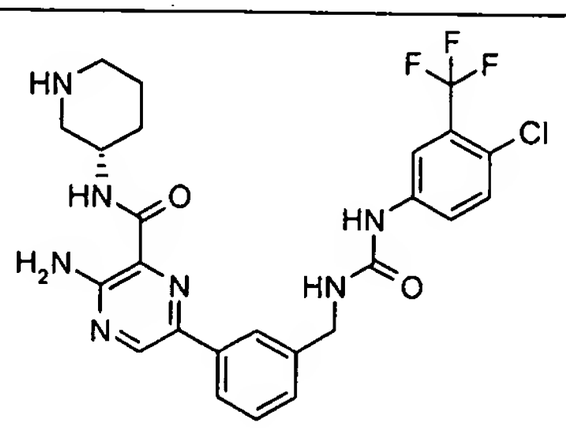
Entry	Name	Structure
178	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[6-(hydroxymethyl)pyridin-3-yl]phenyl}methyl)urea	
179	N-{[3-(6-acetylpyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
180	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(6-cyanopyridin-3-yl)phenyl]methyl}urea	
181	1,1-dimethylethyl (3S)-3-({[3-amino-6-(3-{[({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)pyrazin-2-yl]carbonyl}amino)piperidine-1-carboxylate	
182	3-amino-6-(3-{[({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)-N-[(3S)-piperidin-3-yl]pyrazine-2-carboxamide	

Table 3

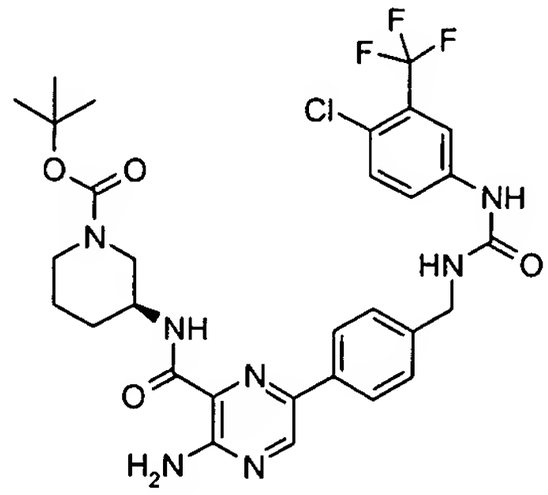
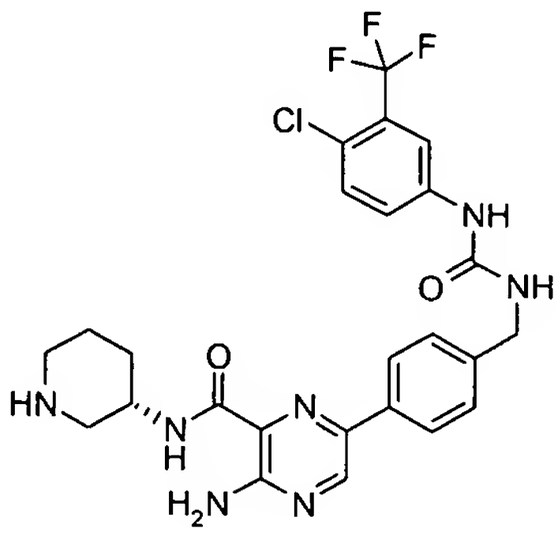
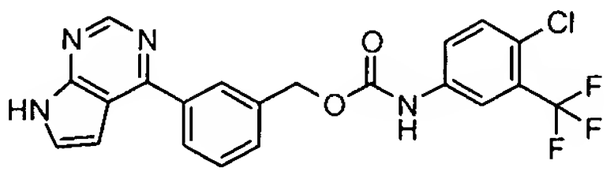
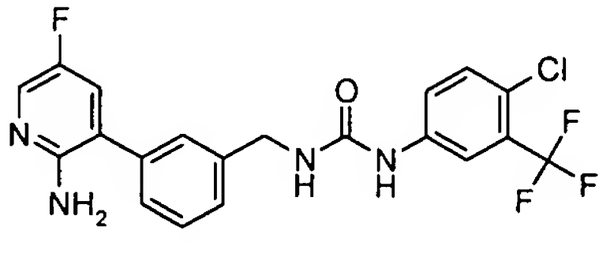
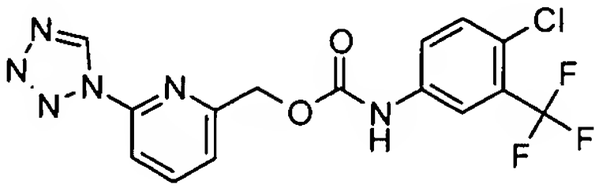
Entry	Name	Structure
183	1,1-dimethylethyl (3S)-3-({[3-amino-6-(4-{[({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)pyrazin-2-yl]carbonyl}amino)piperidine-1-carboxylate	
184	3-amino-6-(4-{[({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)-N-[(3S)-piperidin-3-yl]pyrazine-2-carboxamide	
185	[3-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
186	N-{[3-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
187	[6-(1H-tetrazol-1-yl)pyridin-2-yl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

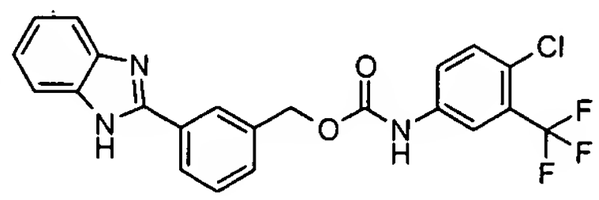
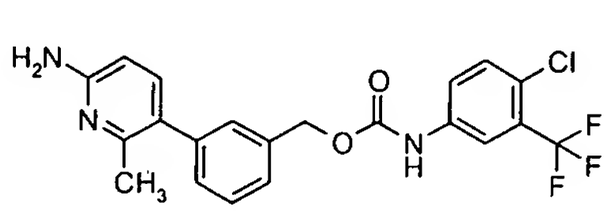
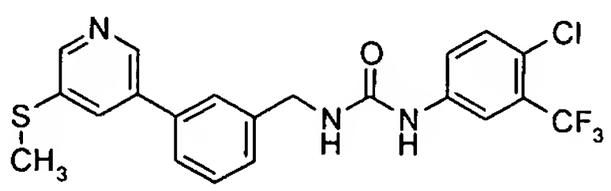
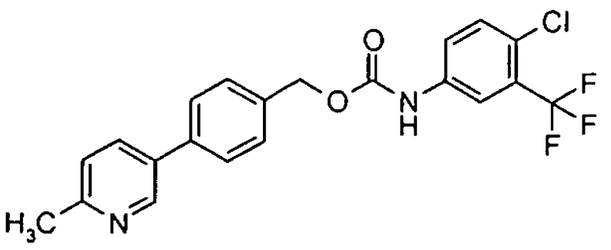
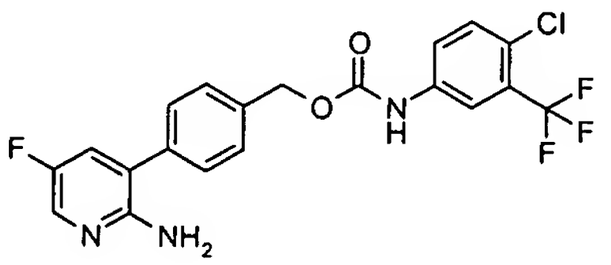
Entry	Name	Structure
188	[3-(1H-benzimidazol-2-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
189	[3-(6-amino-2-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
190	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[5-(methylthio)pyridin-3-yl]phenyl}methyl)urea	
191	[4-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
192	[4-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

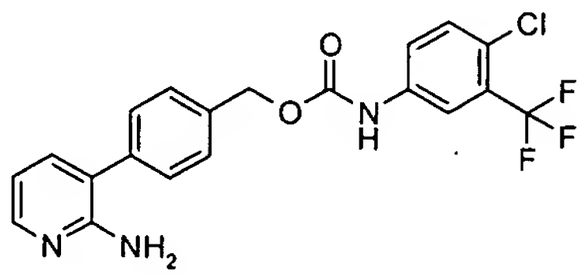
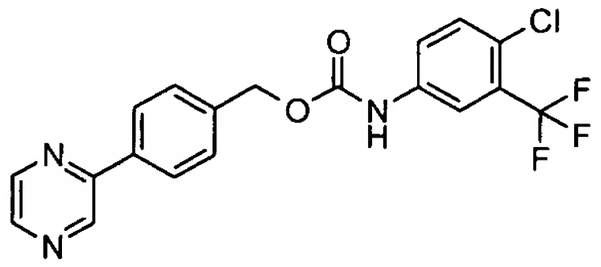
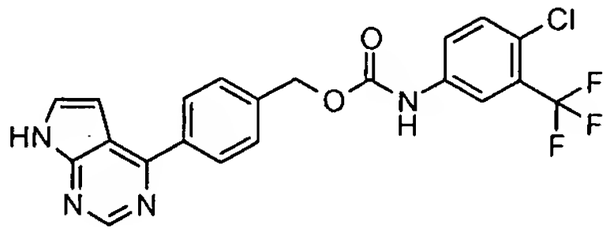
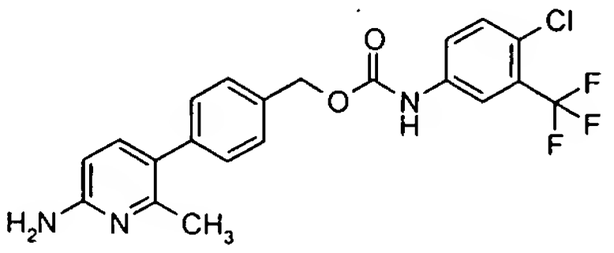
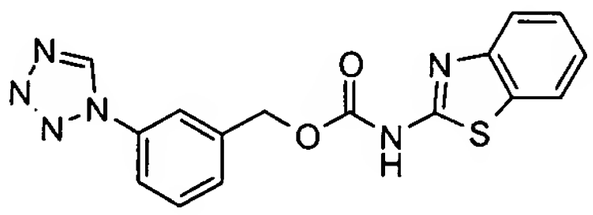
Entry	Name	Structure
193	[4-(2-aminopyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
194	(4-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
195	[4-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
196	[4-(6-amino-2-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
197	[3-(1H-tetrazol-1-yl)phenyl]methyl 1,3-benzothiazol-2-ylcarbamate	

Table 3

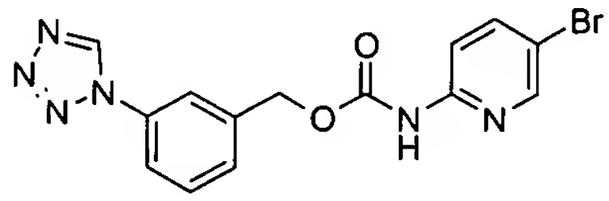
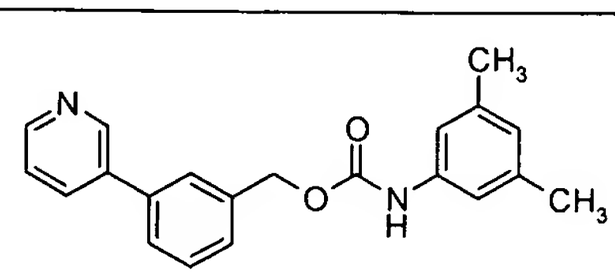
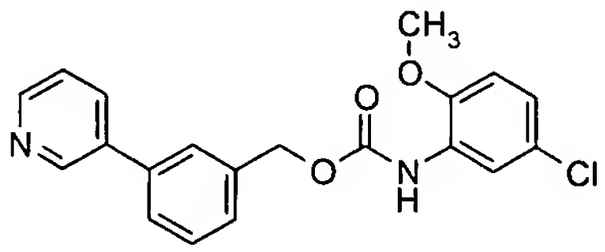
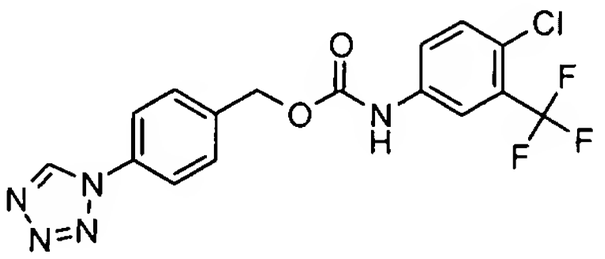
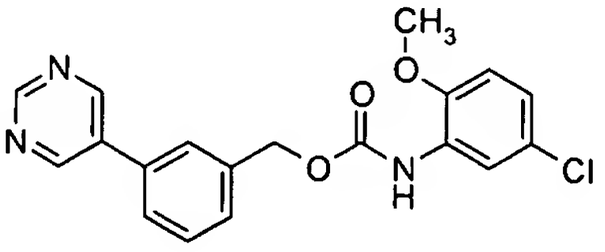
Entry	Name	Structure
198	[3-(1H-tetrazol-1-yl)phenyl]methyl (5-bromopyridin-2-yl)carbamate	
199	(3-pyridin-3-ylphenyl)methyl (3,5-dimethylphenyl)carbamate	
200	(3-pyridin-3-ylphenyl)methyl [5-chloro-2-(methyloxy)phenyl]carbamate	
201	[4-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
202	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2-(methyloxy)phenyl]carbamate	

Table 3

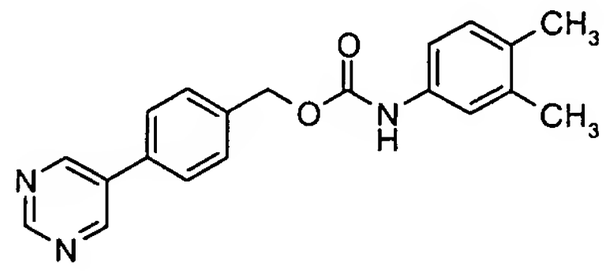
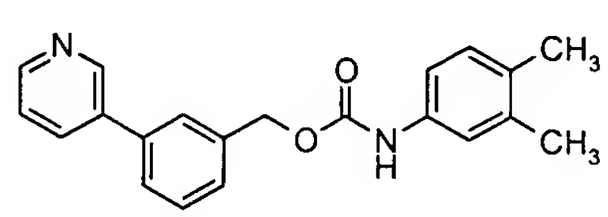
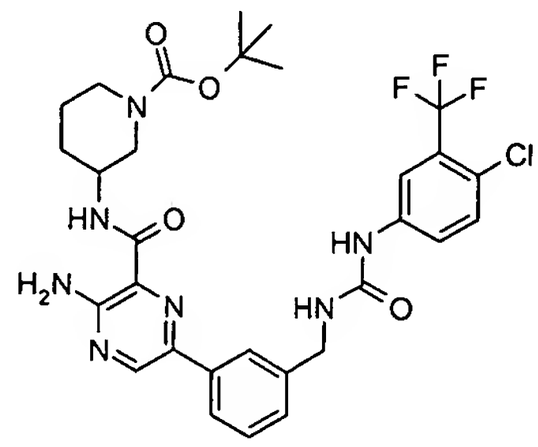
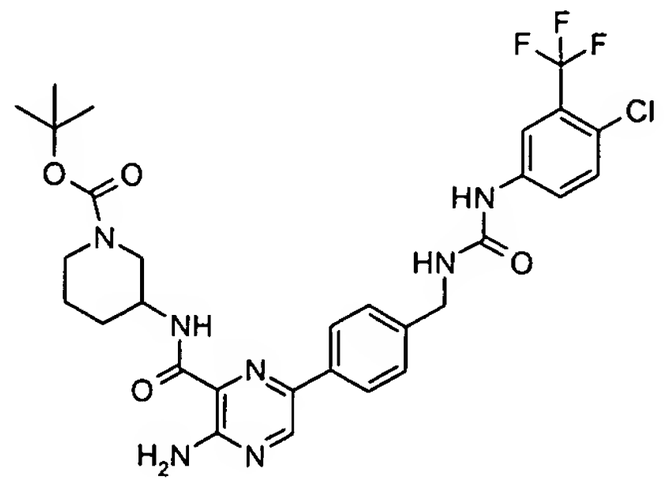
Entry	Name	Structure
203	(4-pyrimidin-5-ylphenyl)methyl (3,4-dimethylphenyl)carbamate	
204	(3-pyridin-3-ylphenyl)methyl (3,4-dimethylphenyl)carbamate	
205	1,1-dimethylethyl 3-({[3-amino-6-(3-{[({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)pyrazin-2-yl]carbonyl}amino)piperidine-1-carboxylate	
206	1,1-dimethylethyl 3-({[3-amino-6-(4-{[({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)pyrazin-2-yl]carbonyl}amino)piperidine-1-carboxylate	

Table 3

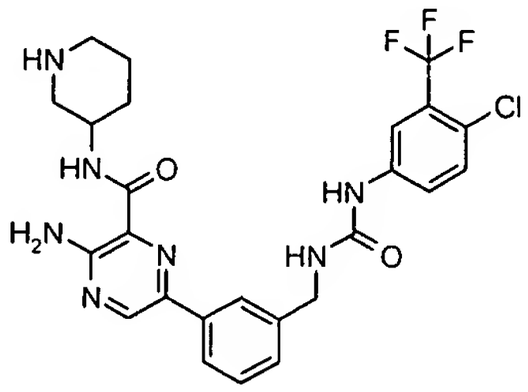
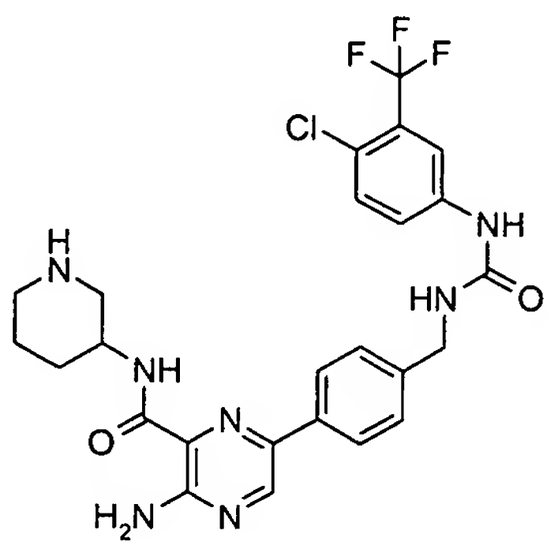
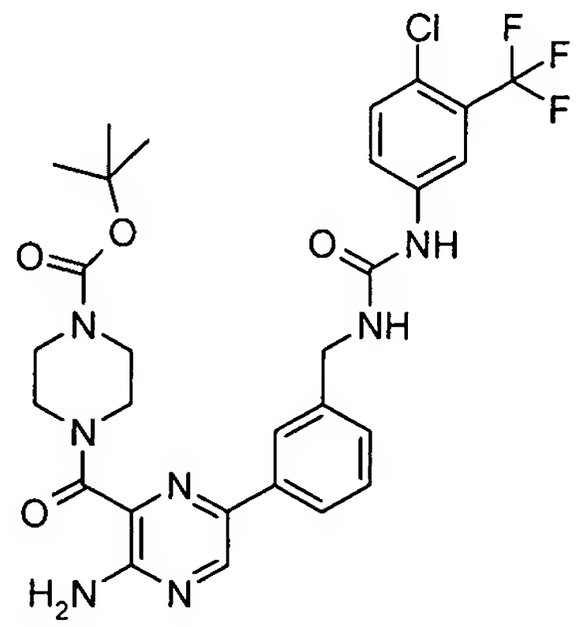
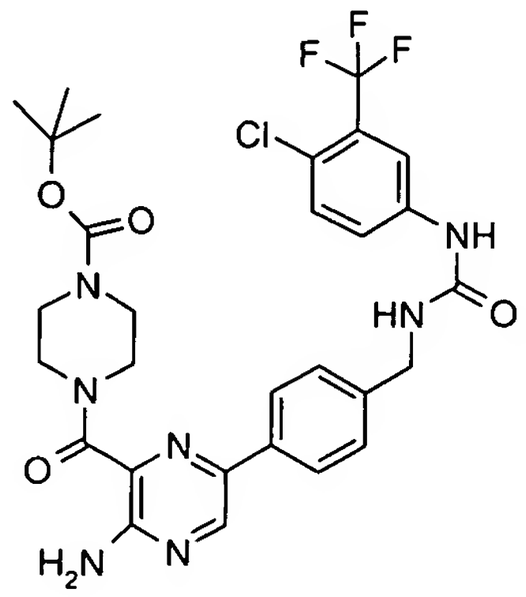
Entry	Name	Structure
207	3-amino-6-(3-{[({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)-N-piperidin-3-ylpyrazine-2-carboxamide	
208	3-amino-6-(4-{[({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)-N-piperidin-3-ylpyrazine-2-carboxamide	
209	1,1-dimethylethyl 4-{[3-amino-6-(3-{[({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)pyrazin-2-yl]carbonyl}piperazine-1-carboxylate	
210	1,1-dimethylethyl 4-{[3-amino-6-(4-{[({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)pyrazin-2-yl]carbonyl}piperazine-1-carboxylate	

Table 3

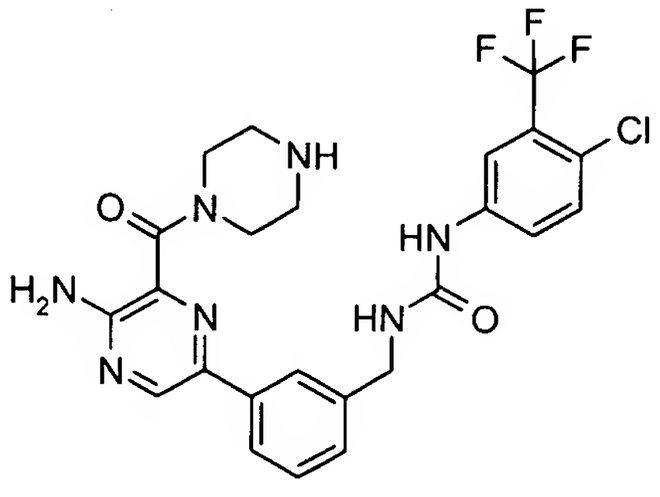
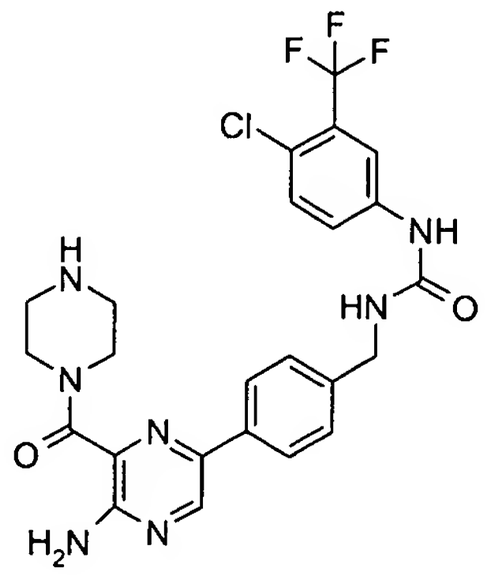
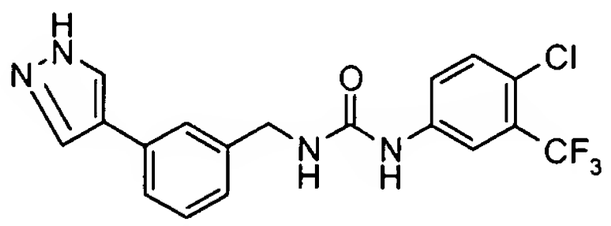
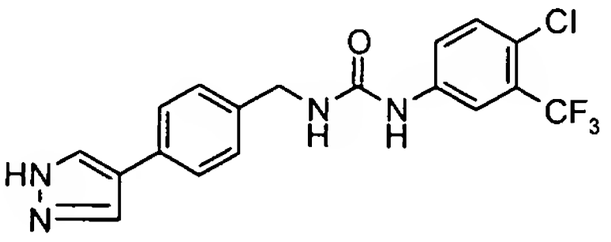
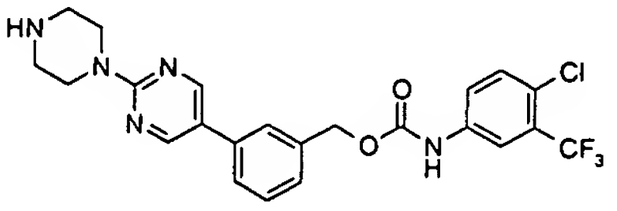
Entry	Name	Structure
211	N-({3-[5-amino-6-(piperazin-1-ylcarbonyl)pyrazin-2-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
212	N-({4-[5-amino-6-(piperazin-1-ylcarbonyl)pyrazin-2-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
213	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(1H-pyrazol-4-yl)phenyl]methyl}urea	
214	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(1H-pyrazol-4-yl)phenyl]methyl}urea	
215	[3-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

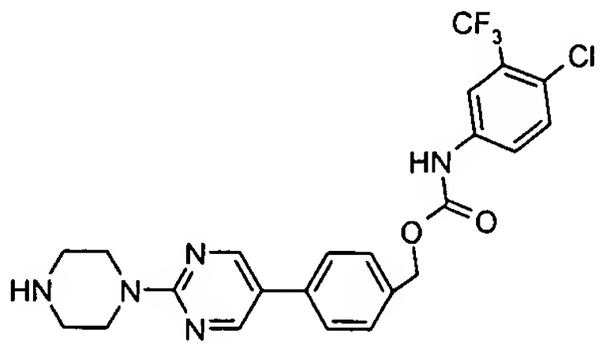
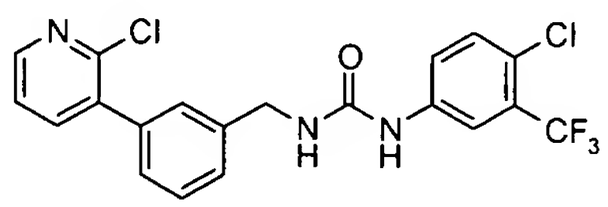
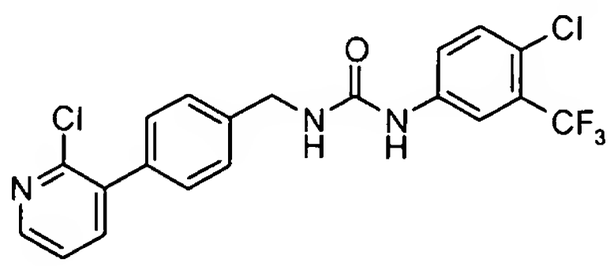
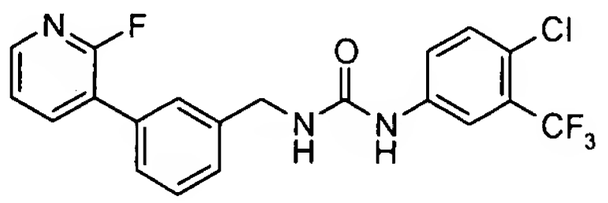
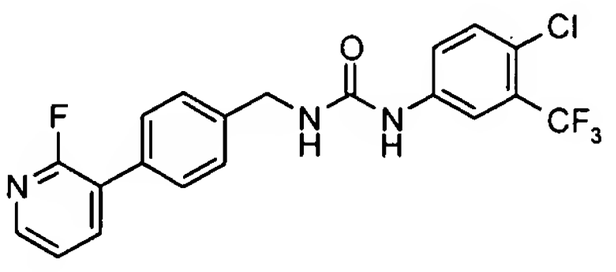
Entry	Name	Structure
216	[4-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
217	N-{[3-(2-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
218	N-{[4-(2-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
219	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(2-fluoropyridin-3-yl)phenyl]methyl}urea	
220	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(2-fluoropyridin-3-yl)phenyl]methyl}urea	

Table 3

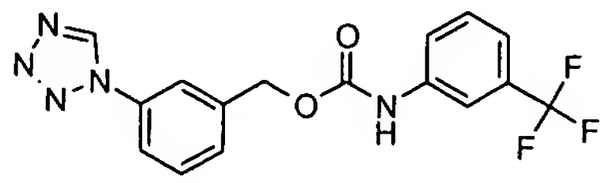
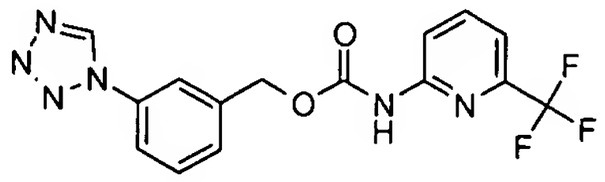
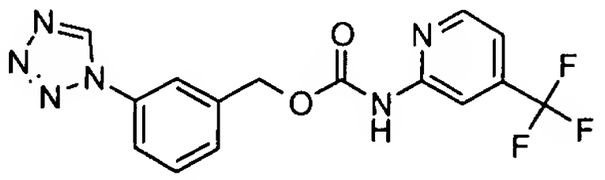
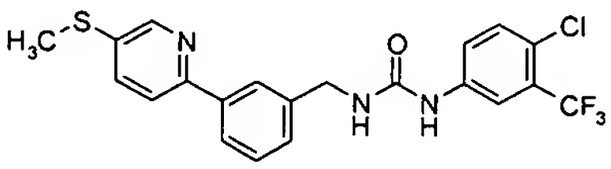
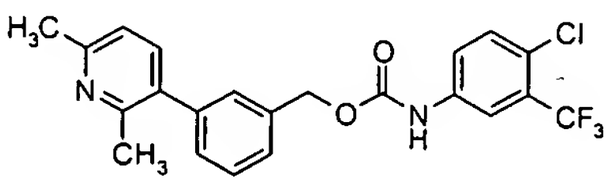
Entry	Name	Structure
221	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-(trifluoromethyl)phenyl]carbamate	
222	[3-(1H-tetrazol-1-yl)phenyl]methyl [6-(trifluoromethyl)pyridin-2-yl]carbamate	
223	[3-(1H-tetrazol-1-yl)phenyl]methyl [4-(trifluoromethyl)pyridin-2-yl]carbamate	
224	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[5-(methylthio)pyridin-2-yl]phenyl}methyl)urea	
225	[3-(2,6-dimethylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

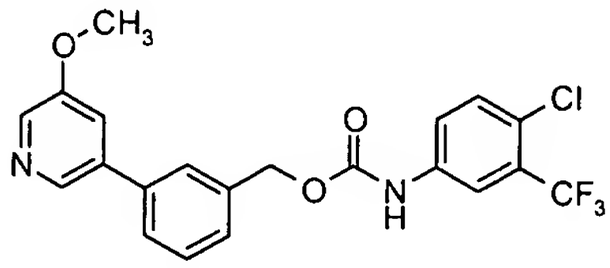
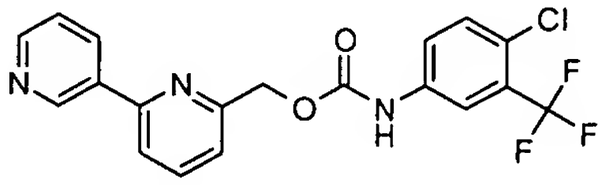
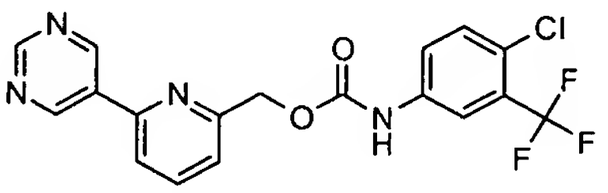
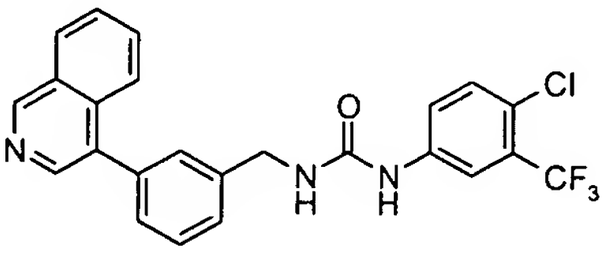
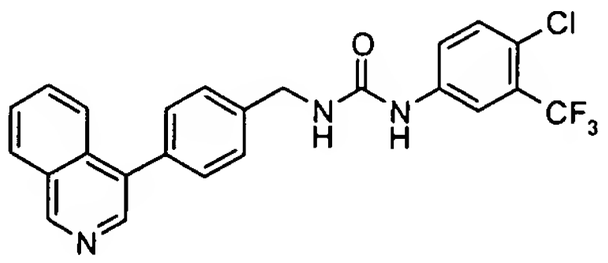
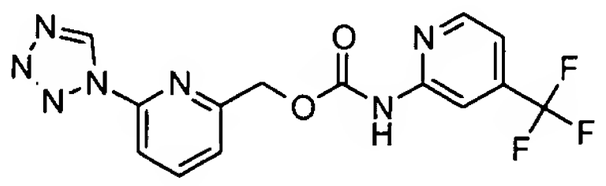
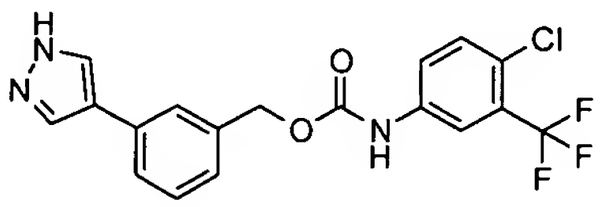
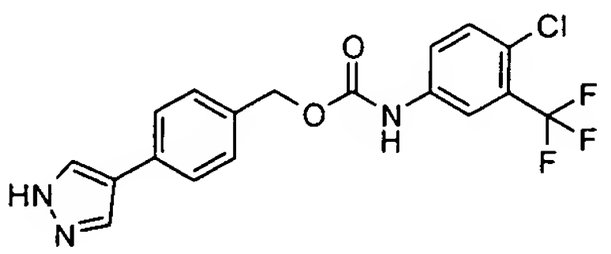
Entry	Name	Structure
226	{3-[5-(methyloxy)pyridin-3-yl]phenyl}methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
227	2,3'-bipyridin-6-ylmethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
228	(6-pyrimidin-5-ylpyridin-2-yl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
229	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-isoquinolin-4-ylphenyl)methyl]urea	
230	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-isoquinolin-4-ylphenyl)methyl]urea	

Table 3

Entry	Name	Structure
231	[6-(1H-tetrazol-1-yl)pyridin-2-yl]methyl [4-(trifluoromethyl)pyridin-2-yl]carbamate	
232	[3-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
233	[4-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

28. (currently amended) A pharmaceutical composition comprising the compound according to ~~any one of claims 1 – 27~~ claim 1 and a pharmaceutically acceptable carrier.

29. (currently amended) A metabolite of the compound or the pharmaceutical composition according to ~~any one of claims 1 – 28~~claim 1.

30. (currently amended) A method for modulating the *in-vivo* activity of a kinase, the method comprising administering to a subject an effective amount of the compound according to ~~any of claims 1 – 27~~claim 1 or a compound selected from N-naphthalen-1-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(phenyloxy)phenyl] -2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,4-dimethylphenyl)-2-{[3-

(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-dimethyl-phenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl] oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl) phenyl]oxy} acetamide, N-(4-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-diethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(ethyloxy) phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(dimethylamino)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-chloro-3-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-bromophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoro-methyl)phenyl] acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl) phenyl] acetamide, methyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, ethyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, 3-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acetamide, N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl] oxy} acetamide, N-(4-aminophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and N-(4-acetylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide.

31. (original) The method according to claim 30, wherein the kinase is c-Kit.

32. (original) The method according to claim 31, wherein modulating the *in vivo* activity of c-Kit comprises inhibition of c-Kit.

33. (currently amended) A method of treating diseases or disorders associated with uncontrolled, abnormal, and/or unwanted cellular activities, the method comprising

administering, to a mammal in need thereof, a therapeutically effective amount of the compound or the pharmaceutical composition as described in ~~any one of claims 1 – 28~~claim 1 or a compound, or a pharmaceutical composition comprising said compound, selected from N-naphthalen-1-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(phenyloxy)phenyl] -2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dimethyl-phenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-dimethyl-phenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-(2,4,6-trimethylphenyl) acetamide, N-(2-ethyl-phenyl)-2-{[3-(1H-tetrazol-1-yl) phenyl]oxy} acetamide, N-(4-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-diethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(ethyloxy) phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2,4-bis(methyl-oxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(dimethylamino)-phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-chloro-3-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-bromophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoro-methyl)phenyl] acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl) phenyl] acetamide, methyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, ethyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, 3-[({[3-(1H-tetrazol-1-yl)phenyl]oxy} acetyl) amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acetamide, N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl] oxy}

acetamide, N-(4-aminophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and N-(4-acetylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide.

34. (currently amended) A method of screening for modulators of c-Kit, the method comprising combining the compound according to ~~any one of claims 1 – 27~~ claim 1 or a compound selected from N-naphthalen-1-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(phenyloxy)phenyl] -2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-dimethyl-phenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl] oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl) phenyl]oxy} acetamide, N-(4-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-diethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(ethyloxy) phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(dimethylamino)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-chloro-3-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-bromophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoro-methyl)phenyl] acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl) phenyl] acetamide, methyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, ethyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, 3-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acetamide, N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl] oxy}

acetamide, N-(4-aminophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and N-(4-acetylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and at least one candidate agent and determining the effect of the candidate agent on c-Kit activity.

35. (currently amended) A method of inhibiting proliferative activity in a cell, the method comprising administering an effective amount of a composition comprising the compound according to claim 1~~any one of claims 1 – 27~~ or a compound selected from N-naphthalen-1-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(phenyloxy)phenyl] -2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-dimethyl-phenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl] oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl) phenyl]oxy} acetamide, N-(4-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-diethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(ethyloxy) phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(dimethylamino)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-chloro-3-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-bromophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoro-methyl)phenyl] acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl) phenyl] acetamide, methyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, ethyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, 3-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{[3-(1H-

tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acetamide, N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl] oxy} acetamide, N-(4-aminophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and N-(4-acetylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, to a cell or a plurality of cells.